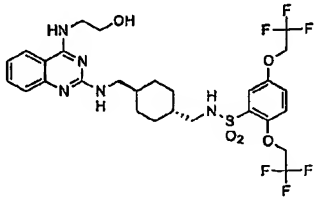
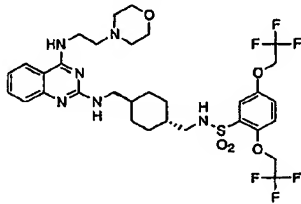
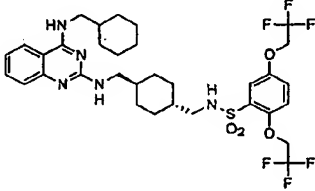
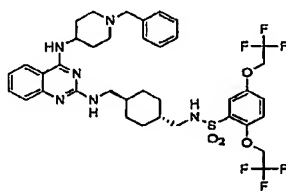
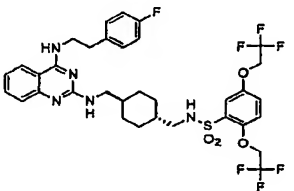
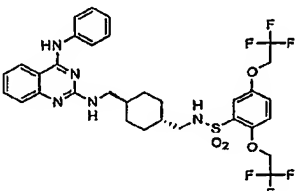
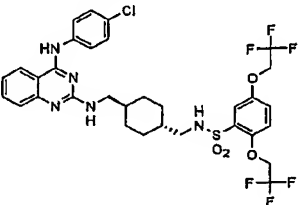
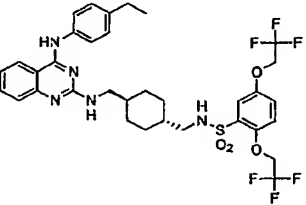
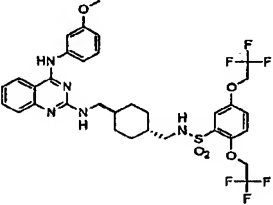
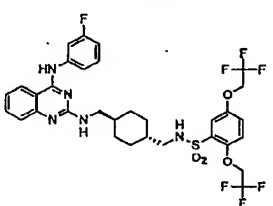
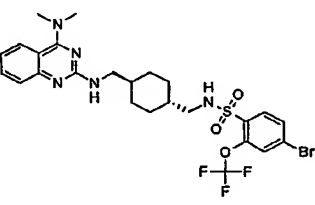
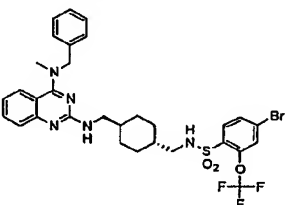
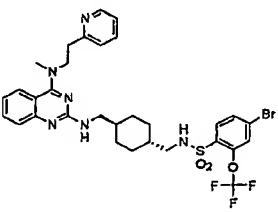
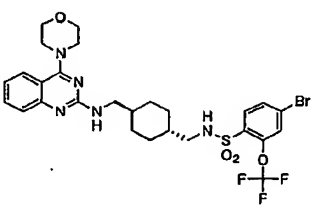
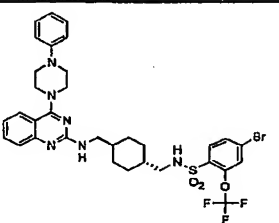
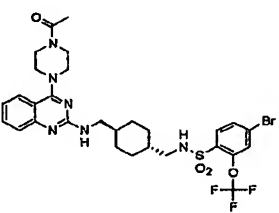
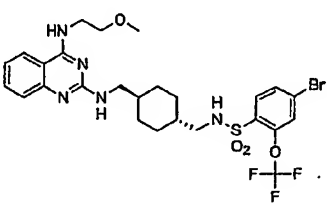
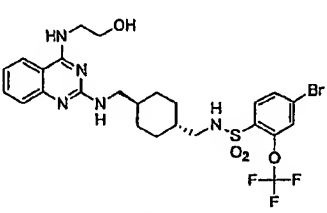
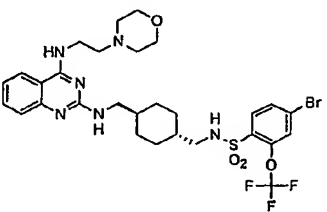
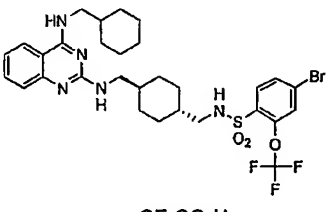
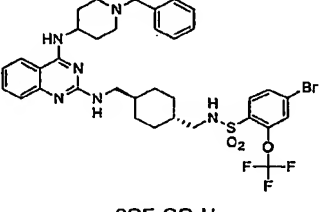
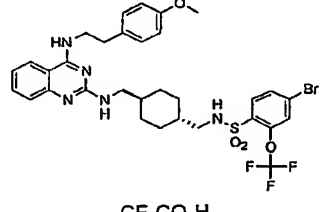
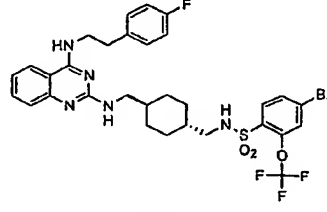
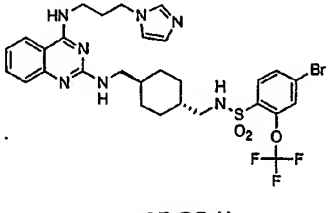
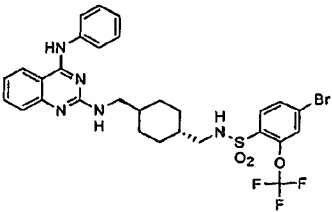
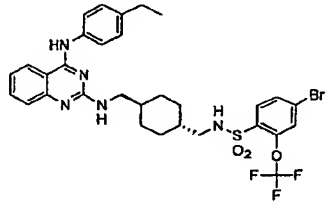
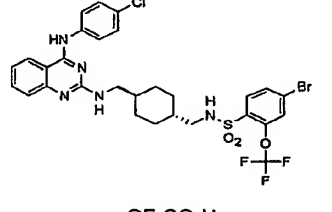
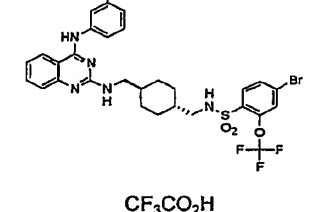
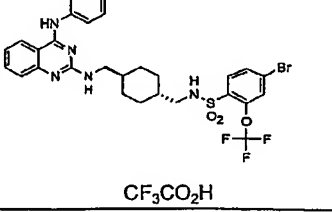
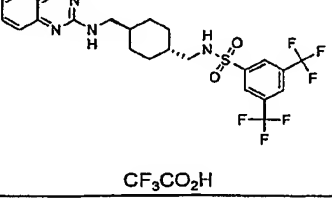


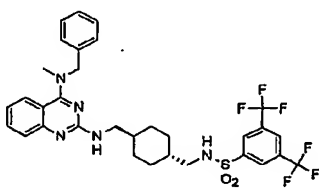
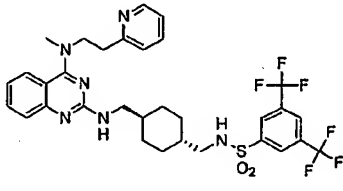
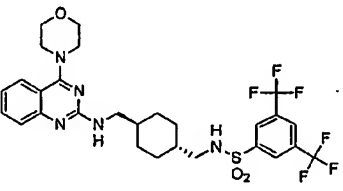
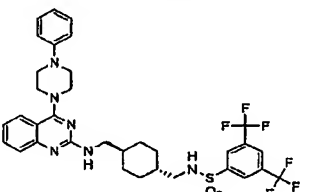
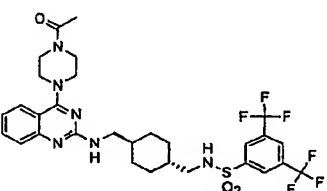
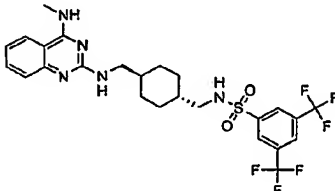
Example No.	Structure	ESI-MS	Retention Time (min)
2423	 <chem>CC1=NC2=C(N1)N=CN=C2C(=N)NCC3CCCCC3NS(=O)(=O)c4cc(OC(F)(F)F)cc(OC(F)(F)F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	666.0 (M + H)	3.86
2424	 <chem>C1CCN(C1)CC2=NC3=CC=CC=C3N=C2C(=N)NCC4CCCCC4NS(=O)(=O)c5cc(OC(F)(F)F)cc(OC(F)(F)F)c5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	735.4 (M + H)	3.50
2425	 <chem>CC1=NC2=CC=CC=C2N=C1C(=N)NCC3CCCCC3NS(=O)(=O)c4cc(OC(F)(F)F)cc(OC(F)(F)F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	718.4 (M + H)	4.64
2426	 <chem>C1CCN(C1Cc2ccccc2)CC3=NC4=CC=CC=C4N=C3C(=N)NCC5CCCCC5NS(=O)(=O)c6cc(OC(F)(F)F)cc(OC(F)(F)F)c6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	795.6 (M + H)	3.70
2427	 <chem>Fc1ccc(cc1)CC2=NC3=CC=CC=C3N=C2C(=N)NCC4CCCCC4NS(=O)(=O)c5cc(OC(F)(F)F)cc(OC(F)(F)F)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	744.2 (M + H)	4.43
2428	 <chem>c1ccc(cc1)CC2=NC3=CC=CC=C3N=C2C(=N)NCC4CCCCC4NS(=O)(=O)c5cc(OC(F)(F)F)cc(OC(F)(F)F)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	698.0 (M + H)	4.26

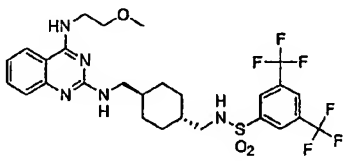
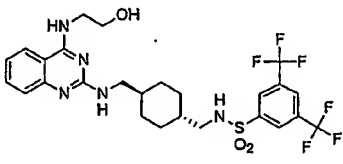
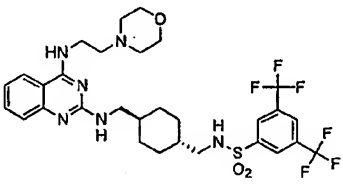
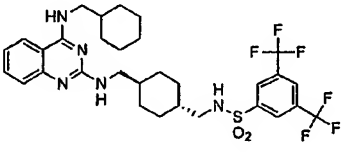
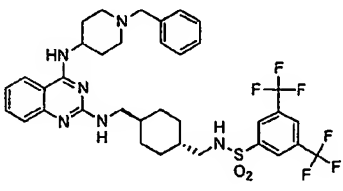
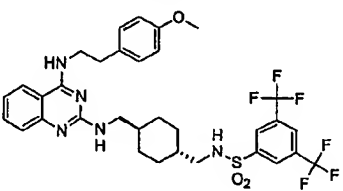
Example No.	Structure	ESI-MS	Retention Time (min)
2429	 <chem>Clc1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	732.4 (M + H)	4.37
2430	 <chem>CCc1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	726.4 (M + H)	4.52
2431	 <chem>COc1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	728.4 (M + H)	4.36
2432	 <chem>Fc1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	716.4 (M + H)	4.32
2433	 <chem>CN(C)c1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	616.0 (M + H)	4.22
2434	 <chem>Brc1ccc(Nc2nc3ccccc3n2CSC(=O)c4ccc(OC(F)(F)F)cc4OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	692.0 (M + H)	4.57

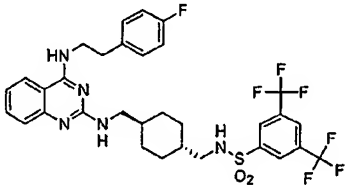
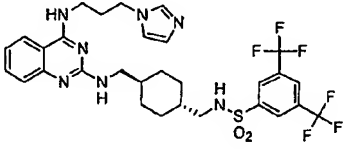
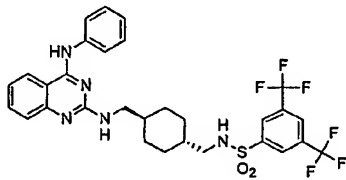
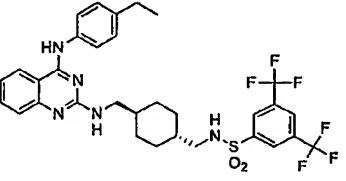
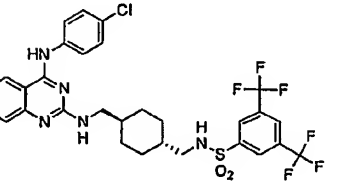
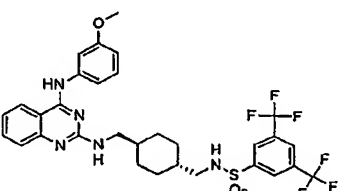
Example No.	Structure	ESI-MS	Retention Time (min)
2435	 2CF ₃ CO ₂ H	707.2 (M + H)	3.64
2436	 CF ₃ CO ₂ H	658.2 (M + H)	4.15
2437	 CF ₃ CO ₂ H	733.2 (M + H)	4.68
2438	 CF ₃ CO ₂ H	699.2 (M + H)	3.88
2439	 CF ₃ CO ₂ H	646.4 (M + H)	4.08
2440	 CF ₃ CO ₂ H	632.4 (M + H)	3.86

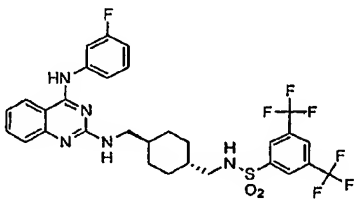
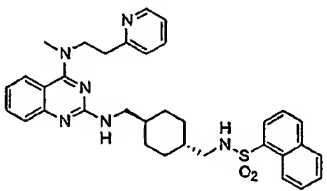
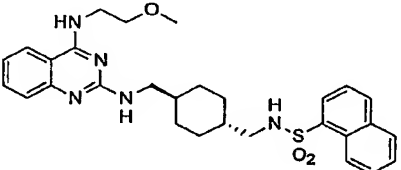
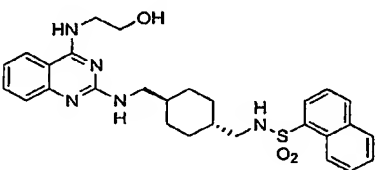
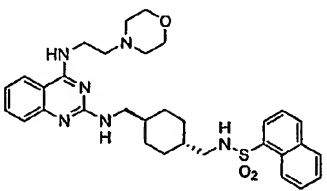
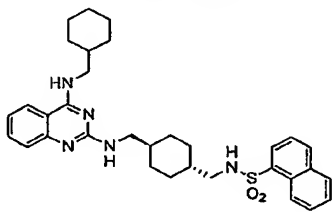
Example No.	Structure	ESI-MS	Retention Time (min)
2441	 $2\text{CF}_3\text{CO}_2\text{H}$	701.4 (M + H)	3.51
2442	 $\text{CF}_3\text{CO}_2\text{H}$	684.2 (M + H)	4.75
2443	 $2\text{CF}_3\text{CO}_2\text{H}$	761.2 (M + H)	3.74
2444	 $\text{CF}_3\text{CO}_2\text{H}$	722.2 (M + H)	4.59
2445	 $\text{CF}_3\text{CO}_2\text{H}$	710.2 (M + H)	4.60
2446	 $2\text{CF}_3\text{CO}_2\text{H}$	696.2 (M + H)	3.53

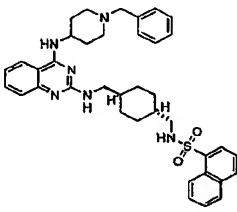
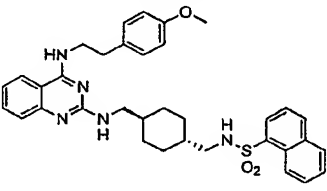
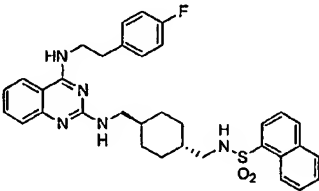
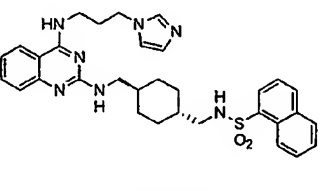
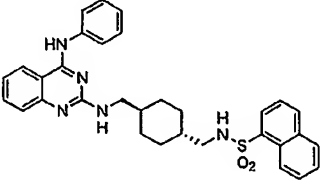
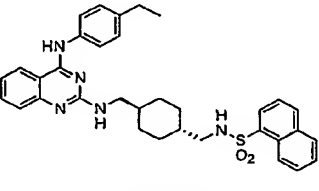
Example No.	Structure	ESI-MS	Retention Time (min)
2447	 <chem>Brc1ccc(cc1)S(=O)(=O)N[C@H]2CCCC[C@H]2CN3C(=N1C(=N2)C(=N1)N3)c4ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	664.2 (M + H)	4.39
2448	 <chem>CCc1ccc(cc1)N2C(=N1C(=N2)C(=N1)N2)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	692.0 (M + H)	4.65
2449	 <chem>Clc1ccc(cc1)N2C(=N1C(=N2)C(=N1)N2)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	698.0 (M + H)	4.59
2450	 <chem>COc1ccc(cc1)N2C(=N1C(=N2)C(=N1)N2)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	694.2 (M + H)	4.42
2451	 <chem>Fc1ccc(cc1)N2C(=N1C(=N2)C(=N1)N2)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	682.2 (M + H)	4.42
2452	 <chem>CN(C)c1nc2ccccc2n1[C@H]3CCCC[C@H]3NS(=O)(=O)c4cc(C(F)(F)F)ccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	590.2 (M + H)	4.28

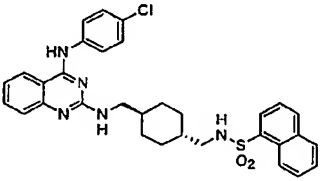
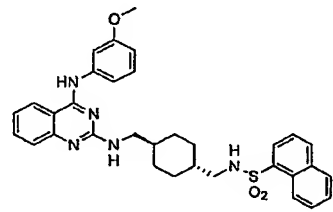
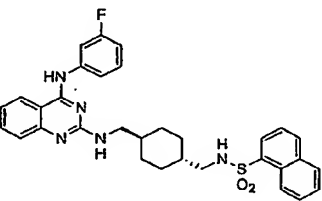
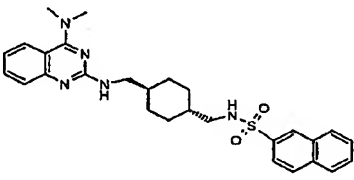
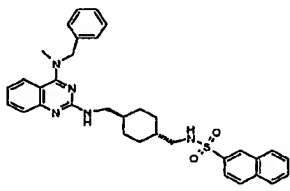
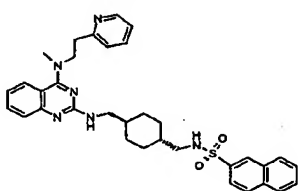
Example No.	Structure	ESI-MS	Retention Time (min)
2453	 $\text{CF}_3\text{CO}_2\text{H}$	666.2 (M + H)	4.61
2454	 $2\text{CF}_3\text{CO}_2\text{H}$	681.2 (M + H)	3.72
2455	 $\text{CF}_3\text{CO}_2\text{H}$	632.4 (M + H)	4.21
2456	 $2\text{CF}_3\text{CO}_2\text{H}$	707.2 (M + H)	4.70
2457	 $\text{CF}_3\text{CO}_2\text{H}$	673.2 (M + H)	3.94
2458	 $\text{CF}_3\text{CO}_2\text{H}$	576.2 (M + H)	4.16

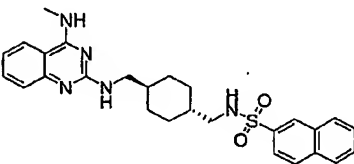
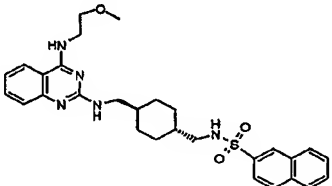
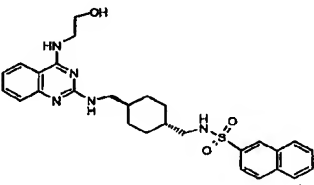
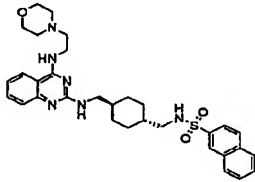
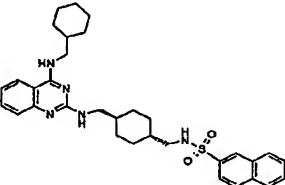
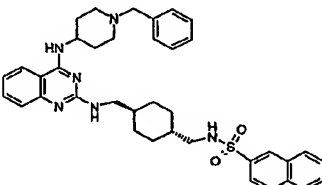
Example No.	Structure	ESI-MS	Retention Time (min)
2459	 <chem>COCCNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3cc(F)c(C(F)(F)F)c(F)c3)c4ccccc14</chem> $\text{CF}_3\text{CO}_2\text{H}$	620.4 (M + H)	4.19
2460	 <chem>OCCNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3cc(F)c(C(F)(F)F)c(F)c3)c4ccccc14</chem> $\text{CF}_3\text{CO}_2\text{H}$	606.6 (M + H)	3.94
2461	 <chem>C1CCN(C1)CCNc2nc3c(ncn3C4CCN(C4)CNS(=O)(=O)c5cc(F)c(C(F)(F)F)c(F)c5)c6ccccc26</chem> $2\text{CF}_3\text{CO}_2\text{H}$	675.4 (M + H)	3.59
2462	 <chem>C1CCN(C1)Cc2nc3c(ncn3C4CCN(C4)CNS(=O)(=O)c5cc(F)c(C(F)(F)F)c(F)c5)c6ccccc26</chem> $\text{CF}_3\text{CO}_2\text{H}$	658.6 (M + H)	4.82
2463	 <chem>C1CCN(C1)Cc2nc3c(ncn3C4CCN(C4)CNS(=O)(=O)c5cc(F)c(C(F)(F)F)c(F)c5)c6ccccc26</chem> $2\text{CF}_3\text{CO}_2\text{H}$	735.4 (M + H)	3.82
2464	 <chem>COc1ccc(cc1)CCNc2nc3c(ncn3C4CCN(C4)CNS(=O)(=O)c5cc(F)c(C(F)(F)F)c(F)c5)c6ccccc26</chem> $\text{CF}_3\text{CO}_2\text{H}$	696.0 (M + H)	4.56

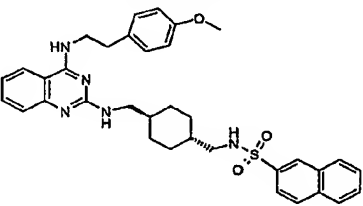
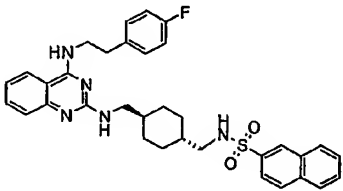
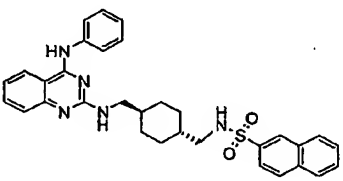
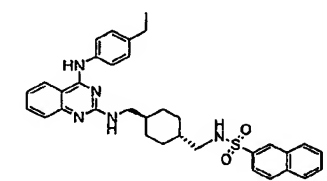
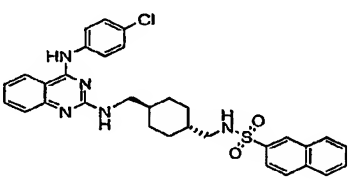
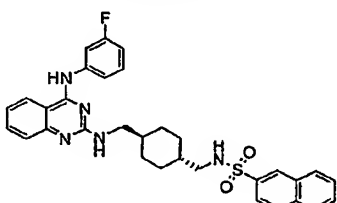
Example No.	Structure	ESI-MS	Retention Time (min)
2465	 <chem>CF3CO2H</chem>	684.4 (M + H)	4.61
2466	 <chem>2CF3CO2H</chem>	670.2 (M + H)	3.56
2467	 <chem>CF3CO2H</chem>	638.2 (M + H)	4.43
2468	 <chem>CF3CO2H</chem>	666.2 (M + H)	4.68
2469	 <chem>CF3CO2H</chem>	672.2 (M + H)	4.60
2470	 <chem>CF3CO2H</chem>	668.2 (M + H)	4.44

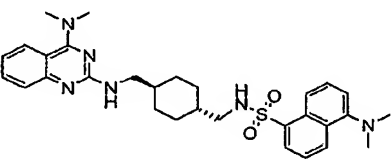
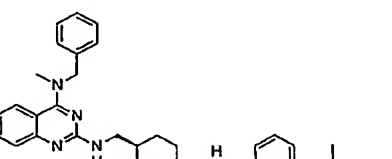
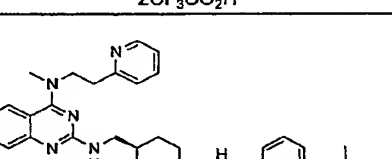
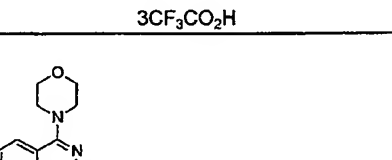
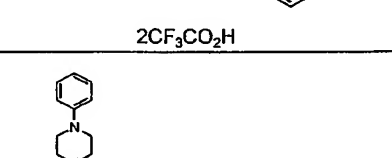
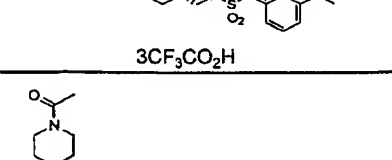
Example No.	Structure	ESI-MS	Retention Time (min)
2471	 $\text{CF}_3\text{CO}_2\text{H}$	656.4 (M + H)	4.47
2472	 $2\text{CF}_3\text{CO}_2\text{H}$	595.4 (M + H)	3.32
2473	 $\text{CF}_3\text{CO}_2\text{H}$	534.0 (M + H)	3.81
2474	 $\text{CF}_3\text{CO}_2\text{H}$	520.4 (M + H)	3.56
2475	 $2\text{CF}_3\text{CO}_2\text{H}$	589.2 (M + H)	3.25
2476	 $\text{CF}_3\text{CO}_2\text{H}$	572.4 (M + H)	4.47

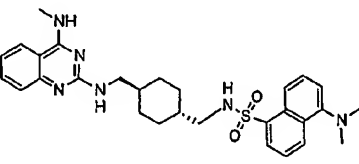
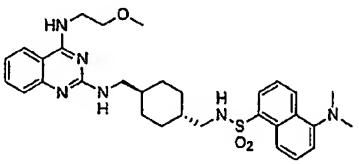
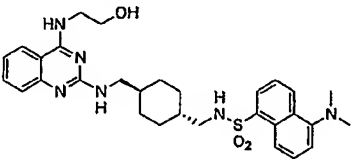
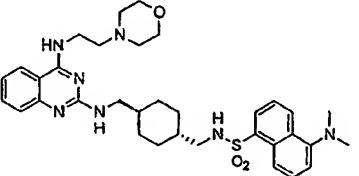
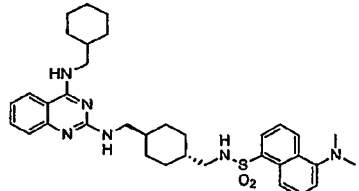
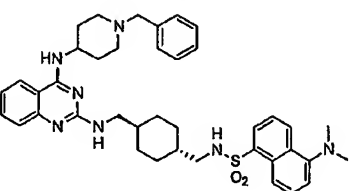
Example No.	Structure	ESI-MS	Retention Time (min)
2477	 <chem>2CF3CO2H</chem>	649.4 (M + H)	3.50
2478	 <chem>CF3CO2H</chem>	610.4 (M + H)	4.26
2479	 <chem>CF3CO2H</chem>	598.2 (M + H)	4.30
2480	 <chem>2CF3CO2H</chem>	584.4 (M + H)	3.29
2481	 <chem>CF3CO2H</chem>	552.6 (M + H)	4.11
2482	 <chem>CF3CO2H</chem>	580.6 (M + H)	4.40

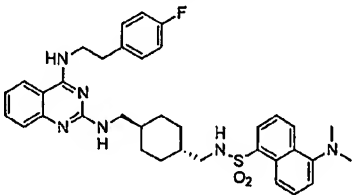
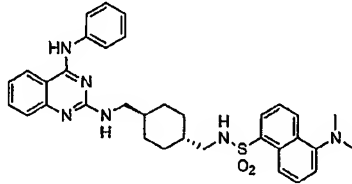
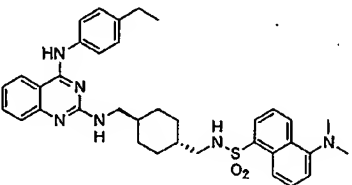
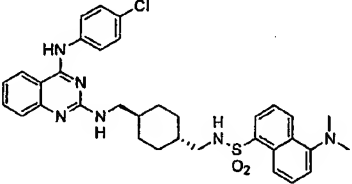
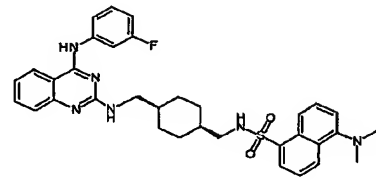
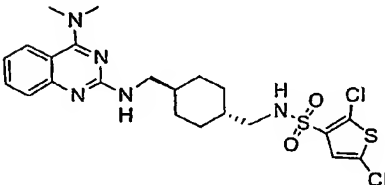
Example No.	Structure	ESI-MS	Retention Time (min)
2483	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNC4CCCCC4NS(=O)(=O)c5ccccc5)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	586.2 (M + H)	4.30
2484	 <chem>COc1ccc(Nc2nc3ccccc3n2CNC4CCCCC4NS(=O)(=O)c5ccccc5)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	582.4 (M + H)	4.14
2485	 <chem>Fc1ccc(Nc2nc3ccccc3n2CNC4CCCCC4NS(=O)(=O)c5ccccc5)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	570.2 (M + H)	4.14
2486	 <chem>CN1C=NC2=C(N1)C=NC2CNC3CCCCC3NS(=O)(=O)c4ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	504.2 (M + H)	3.94
2487	 <chem>CN(C)Cc1ccccc1N2C=NC3=C(N2)C=NC3CNC4CCCCC4NS(=O)(=O)c5ccccc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	580.6 (M + H)	4.34
2488	 <chem>CN(C)Cc1ccccn1N2C=NC3=C(N2)C=NC3CNC4CCCCC4NS(=O)(=O)c5ccccc5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	595.2 (M + H)	3.41

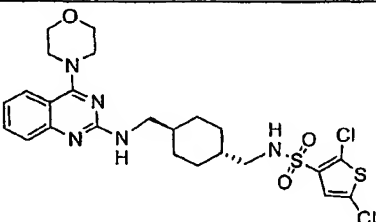
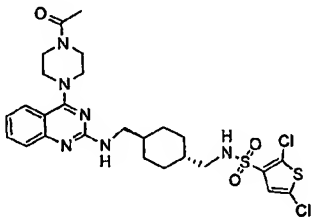
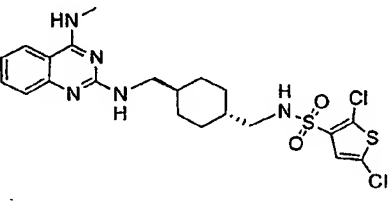
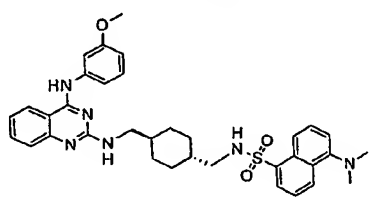
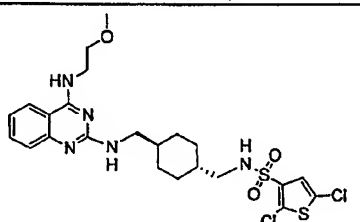
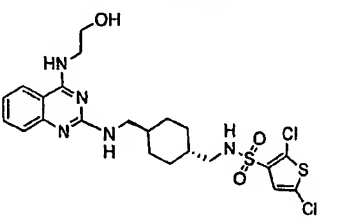
Example No.	Structure	ESI-MS	Retention Time (min)
2489	 <chem>CC1(CCN2C=NC3=CC=CC=C3N=C2N1)CC4(C)CC(C)CC4NS(=O)(=O)c5ccc6ccccc6c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.2 (M + H)	3.84
2490	 <chem>COCCN1C=NC2=CC=CC=C2N=C1N1(C)CC(C)CC1NS(=O)(=O)c3ccc4ccccc4c3</chem> $\text{CF}_3\text{CO}_2\text{H}$	534.2 (M + H)	3.84
2491	 <chem>OCCN1C=NC2=CC=CC=C2N=C1N1(C)CC(C)CC1NS(=O)(=O)c3ccc4ccccc4c3</chem> $\text{CF}_3\text{CO}_2\text{H}$	520.4 (M + H)	3.60
2492	 <chem>C1CCN(C1)CCN2C=NC3=CC=CC=C3N=C2N1(C)CC(C)CC1NS(=O)(=O)c4ccc5ccccc5c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	589.2 (M + H)	3.29
2493	 <chem>C1CCCCC1CN2C=NC3=CC=CC=C3N=C2N1(C)CC(C)CC1NS(=O)(=O)c4ccc5ccccc5c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	572.4 (M + H)	4.51
2494	 <chem>c1ccc(cc1)CN2CCCCC2N3C=NC4=CC=CC=C4N=C3N1(C)CC(C)CC1NS(=O)(=O)c5ccc6ccccc6c5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	649.4 (M + H)	3.52

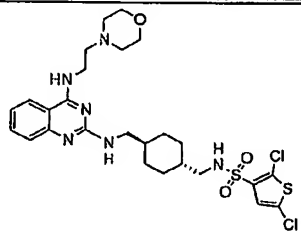
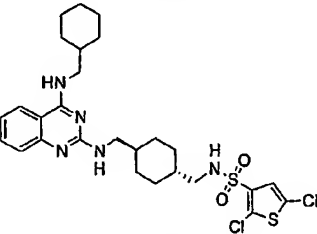
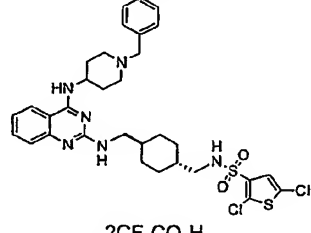
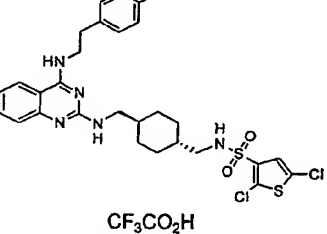
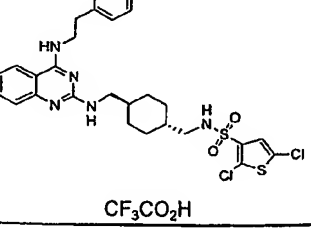
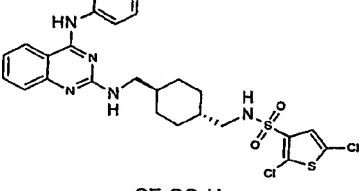
Example No.	Structure	ESI-MS	Retention Time (min)
2495	 CF ₃ CO ₂ H	610.2 (M + H)	4.29
2496	 CF ₃ CO ₂ H	598.2 (M + H)	4.34
2497	 CF ₃ CO ₂ H	552.6 (M + H)	4.13
2498	 CF ₃ CO ₂ H	580.6 (M + H)	4.37
2499	 CF ₃ CO ₂ H	586.2 (M + H)	4.30
2500	 CF ₃ CO ₂ H	570.2 (M + H)	4.18

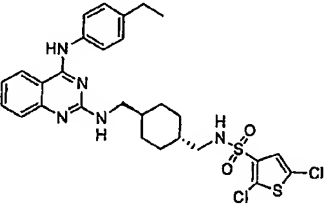
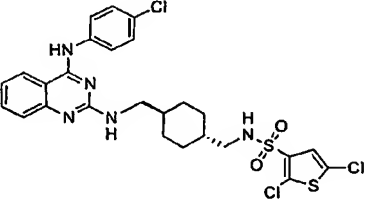
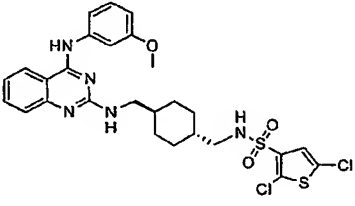
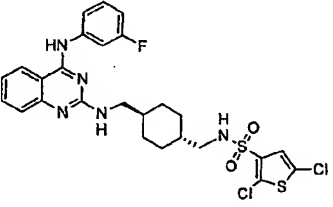
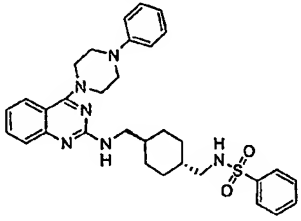
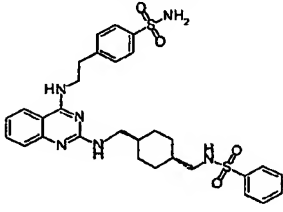
Example No.	Structure	ESI-MS	Retention Time (min)
2501	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	547.4 (M + H)	3.69
2502	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	623.4 (M + H)	4.10
2503	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $3\text{CF}_3\text{CO}_2\text{H}$	638.2 (M + H)	3.20
2504	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	589.2 (M + H)	3.62
2505	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $3\text{CF}_3\text{CO}_2\text{H}$	664.4 (M + H)	4.25
2506	 <chem>CN(C)c1ccc2cc(ccc2c1)S(=O)(=O)N[C@H]3CCCC[C@H]3c4nc5ccccc5n4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	630.4 (M + H)	3.35

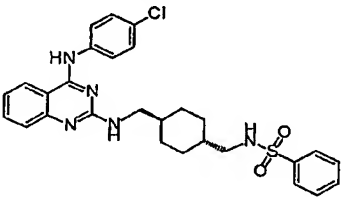
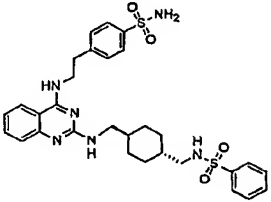
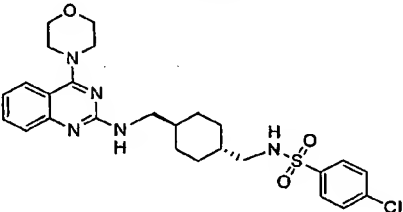
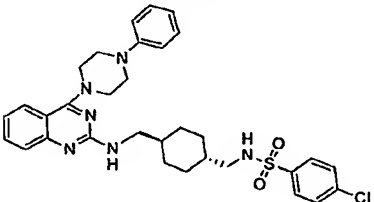
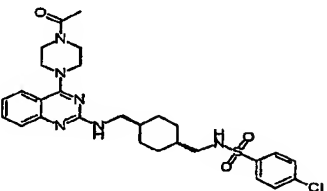
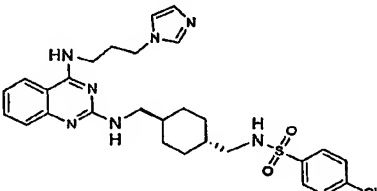
Example No.	Structure	ESI-MS	Retention Time (min)
2507	 $2\text{CF}_3\text{CO}_2\text{H}$	533.2 (M + H)	3.57
2508	 $2\text{CF}_3\text{CO}_2\text{H}$	577.6 (M + H)	3.58
2509	 $2\text{CF}_3\text{CO}_2\text{H}$	563.2 (M + H)	3.28
2510	 $3\text{CF}_3\text{CO}_2\text{H}$	632.6 (M + H)	3.06
2511	 $2\text{CF}_3\text{CO}_2\text{H}$	615.4 (M + H)	4.30
2512	 $3\text{CF}_3\text{CO}_2\text{H}$	692.2 (M + H)	3.38

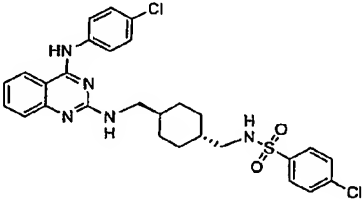
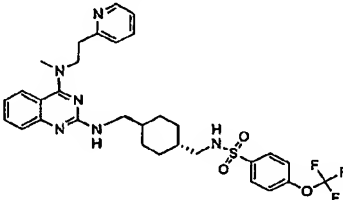
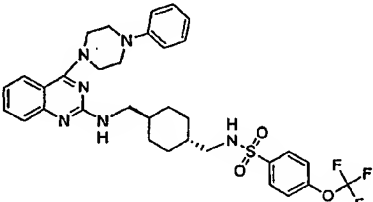
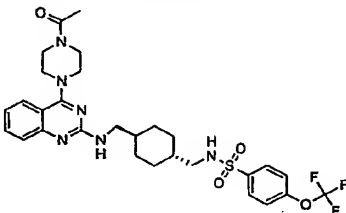
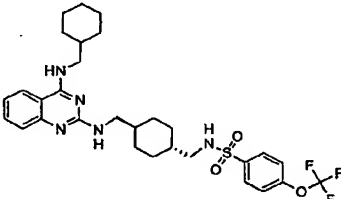
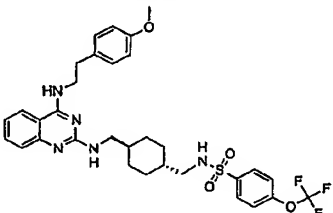
Example No.	Structure	ESI-MS	Retention Time (min)
2513	 $2\text{CF}_3\text{CO}_2\text{H}$	641.4 (M + H)	4.13
2514	 $2\text{CF}_3\text{CO}_2\text{H}$	595.4 (M + H)	3.89
2515	 $2\text{CF}_3\text{CO}_2\text{H}$	623.4 (M + H)	4.20
2516	 $2\text{CF}_3\text{CO}_2\text{H}$	629.2 (M + H)	4.15
2517	 $2\text{CF}_3\text{CO}_2\text{H}$	613.2 (M + H)	4.02
2518	 $\text{CF}_3\text{CO}_2\text{H}$	528.2 (M + H)	4.03

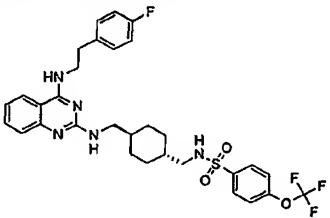
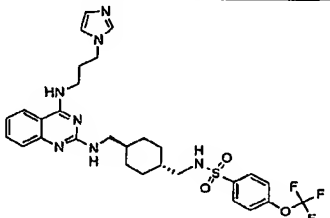
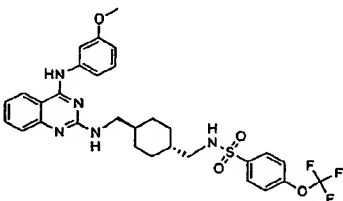
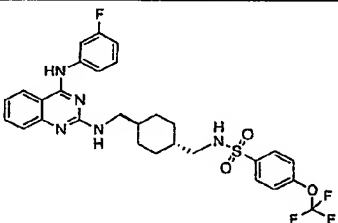
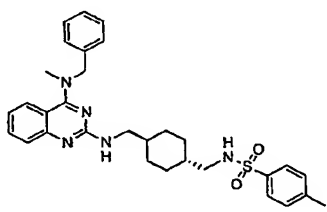
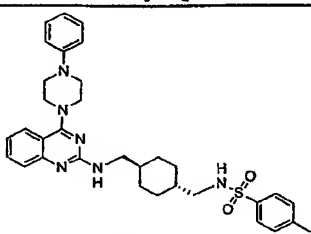
Example No.	Structure	ESI-MS	Retention Time (min)
2519	 <chem>CC1(CCN(C1)c2cc3ccccc3n2)CSC(=O)Nc4cc(Cl)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	570.2 (M + H)	3.96
2520	 <chem>CC1(CCN(C1)c2cc3ccccc3n2)CSC(=O)Nc4cc(Cl)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	611.0 (M + H)	3.69
2521	 <chem>CNc1nc2ccccc2n1CSC(=O)Nc3cc(Cl)cc(Cl)c3</chem> $\text{CF}_3\text{CO}_2\text{H}$	514.2 (M + H)	3.94
2522	 <chem>COc1ccc(Nc2nc3ccccc3n2)cc1CSC(=O)Nc4ccc5c(c4)ccc(NC)cc5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	625.4 (M + H)	3.94
2523	 <chem>COCCNc1nc2ccccc2n1CSC(=O)Nc3cc(Cl)cc(Cl)c3</chem> $\text{CF}_3\text{CO}_2\text{H}$	558.2 (M + H)	3.96
2524	 <chem>OCCNc1nc2ccccc2n1CSC(=O)Nc3cc(Cl)cc(Cl)c3</chem> $\text{CF}_3\text{CO}_2\text{H}$	544.2 (M + H)	3.67

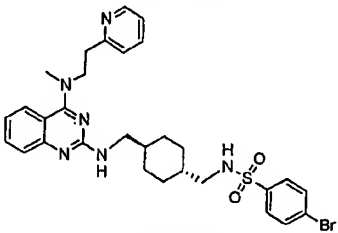
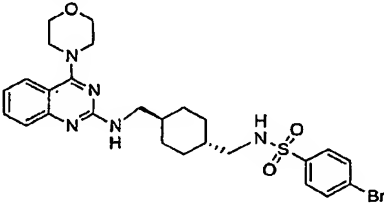
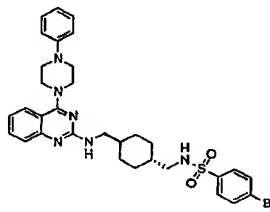
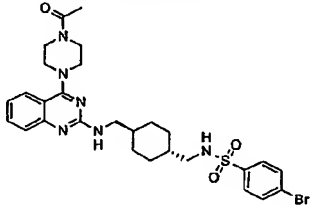
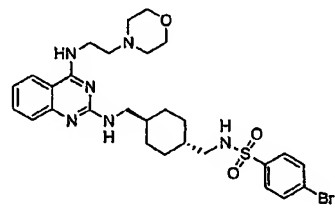
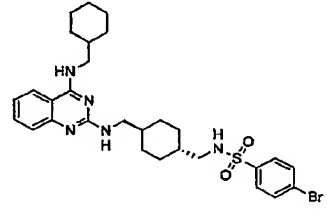
Example No.	Structure	ESI-MS	Retention Time (min)
2525	 $2\text{CF}_3\text{CO}_2\text{H}$	613.2 (M + H)	3.31
2526	 $\text{CF}_3\text{CO}_2\text{H}$	596.2 (M + H)	4.69
2527	 $2\text{CF}_3\text{CO}_2\text{H}$	673.4 (M + H)	3.57
2528	 $\text{CF}_3\text{CO}_2\text{H}$	634.4 (M + H)	4.41
2529	 $\text{CF}_3\text{CO}_2\text{H}$	622.2 (M + H)	4.45
2530	 $\text{CF}_3\text{CO}_2\text{H}$	576 (M + H)	4.25

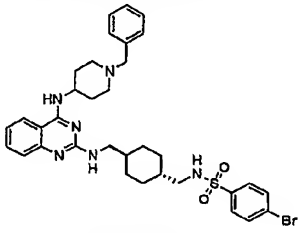
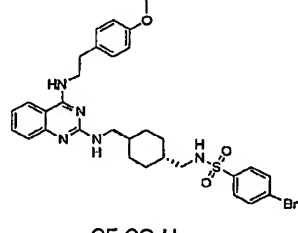
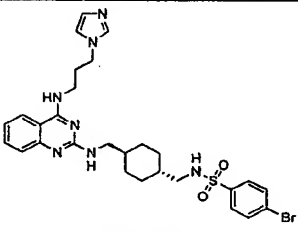
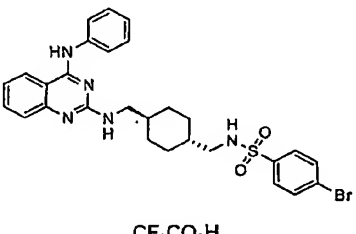
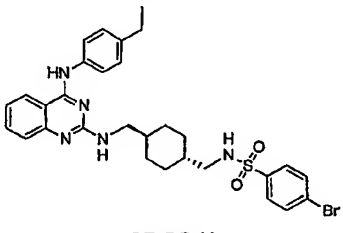
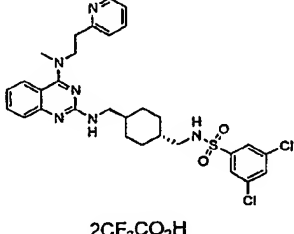
Example No.	Structure	ESI-MS	Retention Time (min)
2531	 <chem>CC1=CC=C(NC2=NC3=CC=CC=C3N=C2NCC4CCCCC4NS(=O)(=O)C5=CC(=C(C=C5)S=C(C=C5)Cl)Cl)C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	604.4 (M + H)	4.52
2532	 <chem>ClC1=CC=C(NC2=NC3=CC=CC=C3N=C2NCC4CCCCC4NS(=O)(=O)C5=CC(=C(C=C5)S=C(C=C5)Cl)Cl)C6=CC=C(C=C6)Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	610.2 (M + H)	4.40
2533	 <chem>COc1ccc(NC2=NC3=CC=CC=C3N=C2NCC4CCCCC4NS(=O)(=O)C5=CC(=C(C=C5)S=C(C=C5)Cl)Cl)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	606.4 (M + H)	4.29
2534	 <chem>Fc1ccc(NC2=NC3=CC=CC=C3N=C2NCC4CCCCC4NS(=O)(=O)C5=CC(=C(C=C5)S=C(C=C5)Cl)Cl)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	594.2 (M + H)	4.27
2535	 <chem>c1ccc(N2CCN(CC2)C3=NC4=CC=CC=C4N=C3NCC5CCCCC5NS(=O)(=O)C6=CC=CC=C6)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	571.8 (M + H)	4.99
2536	 <chem>NC(=O)c1ccc(cc1)SCC2=NC3=CC=CC=C3N=C2NCC4CCCCC4NS(=O)(=O)C5=CC=CC=C5</chem> $\text{CF}_3\text{CO}_2\text{H}$	609.8 (M + H)	4.43

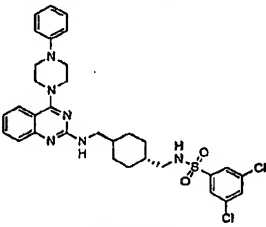
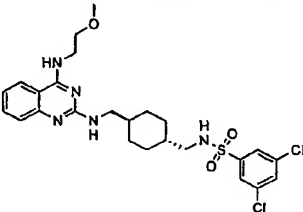
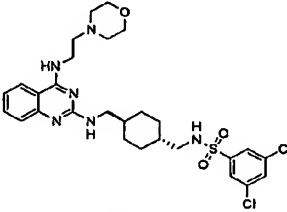
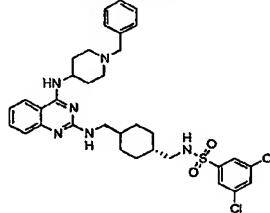
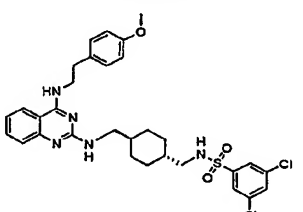
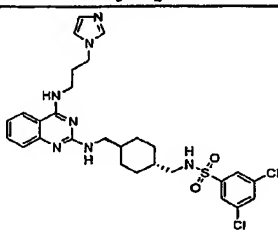
Example No.	Structure	ESI-MS	Retention Time (min)
2537	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	4.86
2538	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	564.6 (M + H)	5.13
2539	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	530.6 (M + H)	4.65
2540	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	605.6 (M + H)	5.21
2541	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	571.6 (M + H)	4.45
2542	 <chem>Clc1ccc(Nc2nc3ccccc3n2CNS(=O)(=O)c4ccc(Cl)cc4)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	568.8 (M + H)	4.09

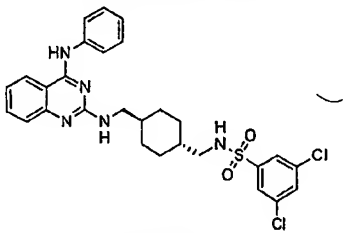
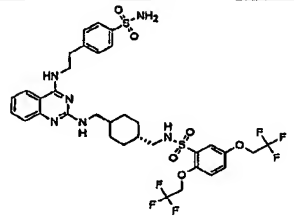
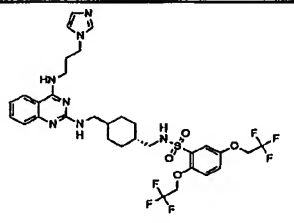
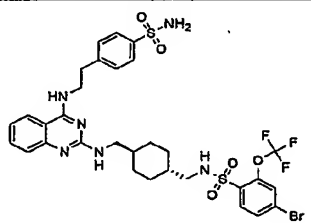
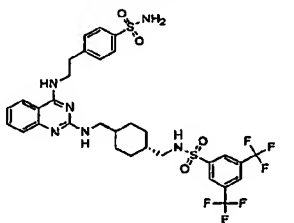
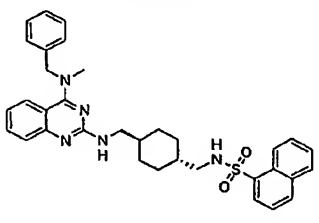
Example No.	Structure	ESI-MS	Retention Time (min)
2543	 CF ₃ CO ₂ H	570.6 (M + H)	5.11
2544	 2CF ₃ CO ₂ H	629.6 (M + H)	4.37
2545	 2CF ₃ CO ₂ H	655.6 (M + H)	5.35
2546	 CF ₃ CO ₂ H	621.8 (M + H)	4.63
2547	 CF ₃ CO ₂ H	606.8 (M + H)	5.45
2548	 CF ₃ CO ₂ H	644.6 (M + H)	5.21

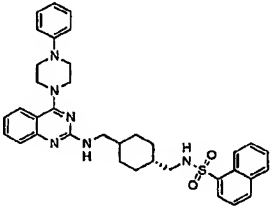
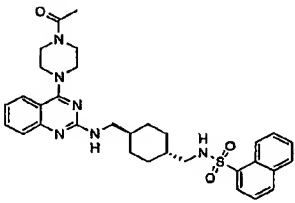
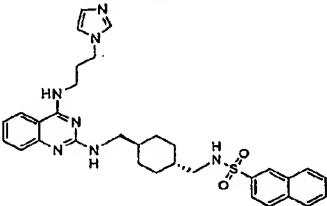
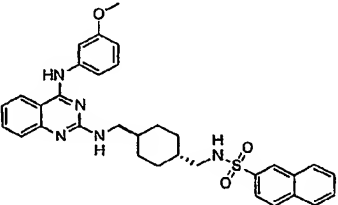
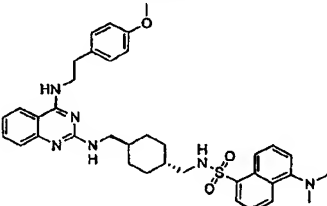
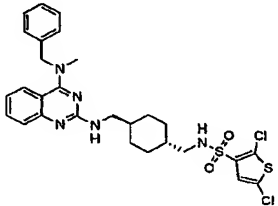
Example No.	Structure	ESI-MS	Retention Time (min)
2549	 $\text{CF}_3\text{CO}_2\text{H}$	632.6 (M + H)	5.25
2550	 $2\text{CF}_3\text{CO}_2\text{H}$	618.6 (M + H)	4.29
2551	 $\text{CF}_3\text{CO}_2\text{H}$	616.6 (M + H)	5.14
2552	 $\text{CF}_3\text{CO}_2\text{H}$	604.6 (M + H)	5.13
2553	 $\text{CF}_3\text{CO}_2\text{H}$	544.6 (M + H)	5.03
2554	 $2\text{CF}_3\text{CO}_2\text{H}$	585.6 (M + H)	5.13

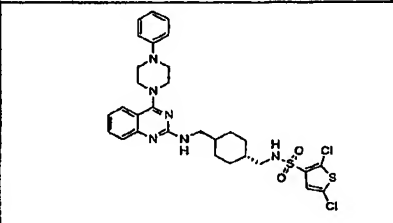
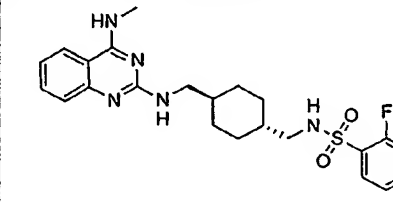
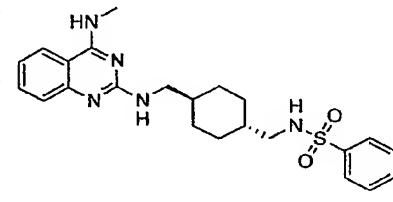
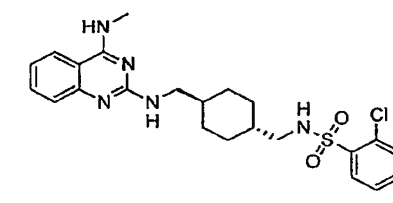
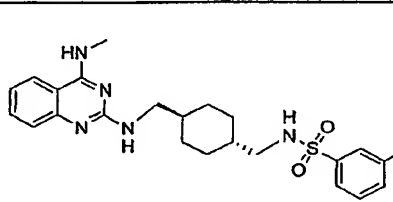
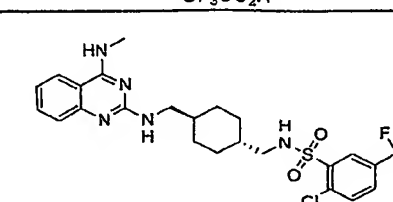
Example No.	Structure	ESI-MS	Retention Time (min)
2555	 $2\text{CF}_3\text{CO}_2\text{H}$	623.6 (M + H)	4.25
2556	 $\text{CF}_3\text{CO}_2\text{H}$	574.6 (M + H)	4.73
2557	 $2\text{CF}_3\text{CO}_2\text{H}$	649.0 (M + H)	5.25
2558	 $\text{CF}_3\text{CO}_2\text{H}$	615.0 (M + H)	4.51
2559	 $2\text{CF}_3\text{CO}_2\text{H}$	617.4 (M + H)	4.15
2560	 $\text{CF}_3\text{CO}_2\text{H}$	600.6 (M + H)	5.37

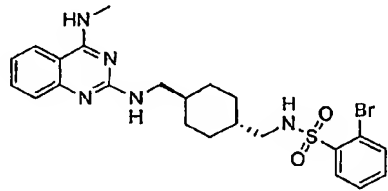
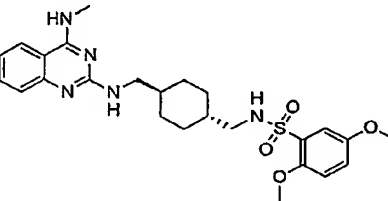
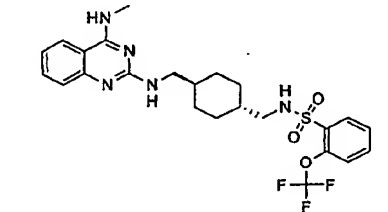
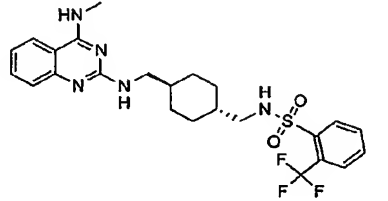
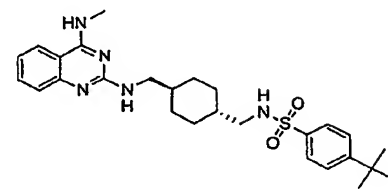
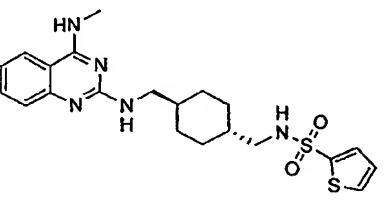
Example No.	Structure	ESI-MS	Retention Time (min)
2561	 $2CF_3CO_2H$	677.0 (M + H)	4.45
2562	 CF_3CO_2H	638.6 (M + H)	5.18
2563	 $2CF_3CO_2H$	612.6 (M + H)	4.16
2564	 CF_3CO_2H	580.0 (M + H)	5.01
2565	 CF_3CO_2H	608.0 (M + H)	5.26
2566	 $2CF_3CO_2H$	613.6 (M + H)	4.44

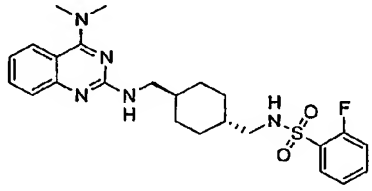
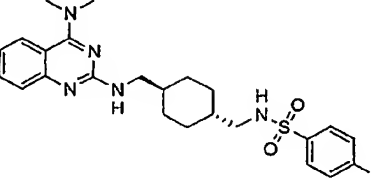
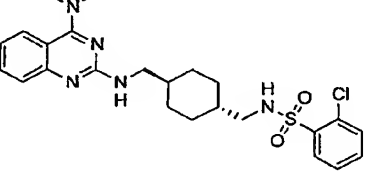
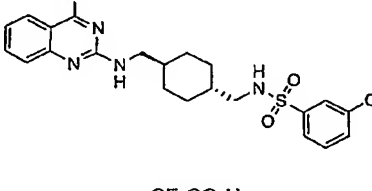
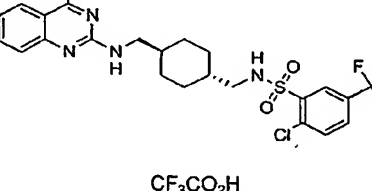
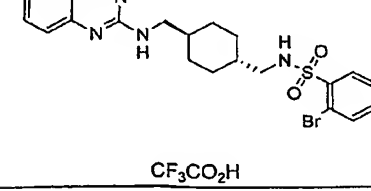
Example No.	Structure	ESI-MS	Retention Time (min)
2567	 2CF ₃ CO ₂ H	639.6 (M + H)	5.48
2568	 CF ₃ CO ₂ H	552.6 (M + H)	4.92
2569	 2CF ₃ CO ₂ H	607.8 (M + H)	4.33
2570	 2CF ₃ CO ₂ H	667.4 (M + H)	4.67
2571	 CF ₃ CO ₂ H	628.6 (M + H)	5.29
2572	 2CF ₃ CO ₂ H	602.6 (M + H)	4.35

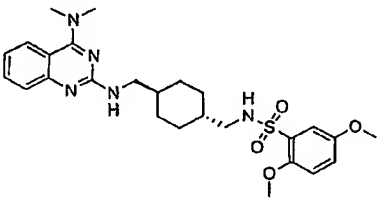
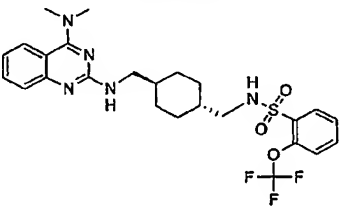
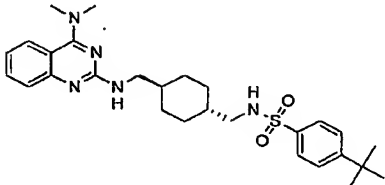
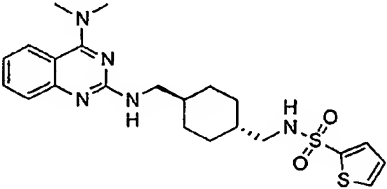
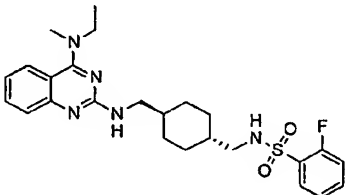
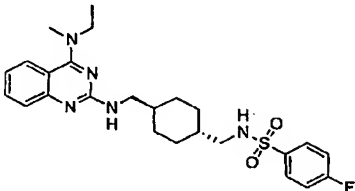
Example No.	Structure	ESI-MS	Retention Time (min)
2573	 <chem>Clc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(Cl)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	570.6 (M + H)	5.23
2574	 <chem>COc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(OC(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	805.4 (M + H)	4.91
2575	 <chem>COc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(OC(F)(F)F)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	730.8 (M + H)	4.47
2576	 <chem>COc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(OC(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	771.6 (M + H)	4.93
2577	 <chem>COc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(OC(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	745.6 (M + H)	5.01
2578	 <chem>COc1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2C3=NC4=CC=CC=C4N(C3)N5=CC=CC=C5N4)c(OC(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	580.8 (M + H)	5.18

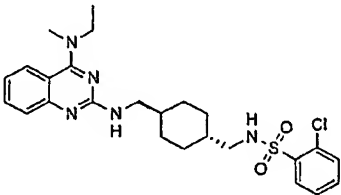
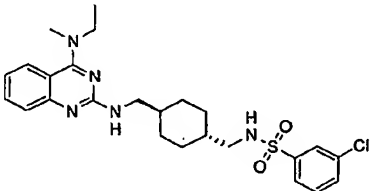
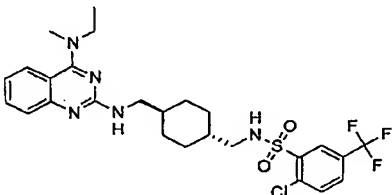
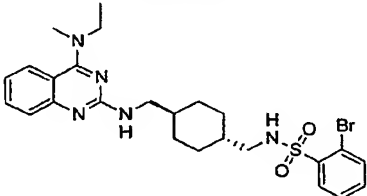
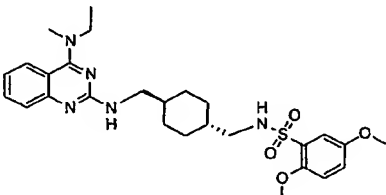
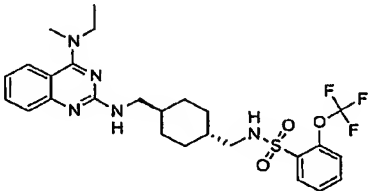
Example No.	Structure	ESI-MS	Retention Time (min)
2579	 $2\text{CF}_3\text{CO}_2\text{H}$	621.8 (M + H)	5.27
2580	 $\text{CF}_3\text{CO}_2\text{H}$	587.6 (M + H)	4.51
2581	 $2\text{CF}_3\text{CO}_2\text{H}$	584.6 (M + H)	4.21
2582	 $\text{CF}_3\text{CO}_2\text{H}$	582.8 (M + H)	5.03
2583	 $\text{CF}_3\text{CO}_2\text{H}$	653.8 (M + H)	4.90
2584	 $\text{CF}_3\text{CO}_2\text{H}$	604.6 (M + H)	5.33

Example No.	Structure	ESI-MS	Retention Time (min)
2585	 $2\text{CF}_3\text{CO}_2\text{H}$	645.6 (M + H)	5.41
2586	 $\text{CF}_3\text{CO}_2\text{H}$	458.6 (M + H)	4.39
2587	 $\text{CF}_3\text{CO}_2\text{H}$	458.6 (M + H)	4.40
2588	 $\text{CF}_3\text{CO}_2\text{H}$	474.6 (M + H)	4.39
2589	 $\text{CF}_3\text{CO}_2\text{H}$	474.6 (M + H)	4.58
2590	 $\text{CF}_3\text{CO}_2\text{H}$	542.6 (M + H)	4.79

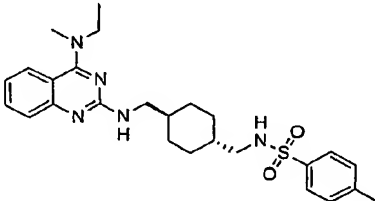
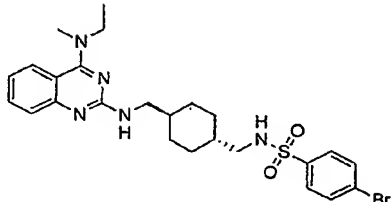
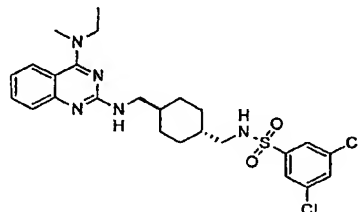
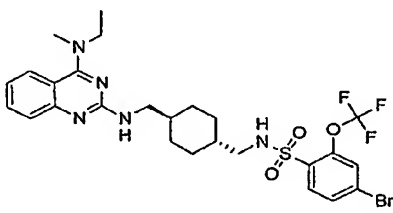
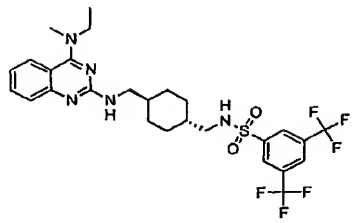
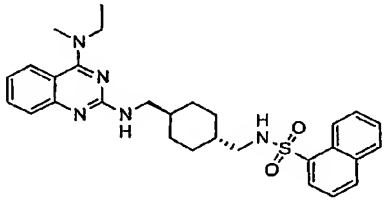
Example No.	Structure	ESI-MS	Retention Time (min)
2591	 <chem>BrC1=CC=C(C=C1)S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.51
2592	 <chem>COc1cc(OC)ccc1S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N</chem> $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.33
2593	 <chem>FC(F)(F)c1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	524.6 (M + H)	4.61
2594	 <chem>Fc1cc(S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N)cc(F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	508.6 (M + H)	4.57
2595	 <chem>CC(C)(C)c1ccc(S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	496.8 (M + H)	4.87
2596	 <chem>c1cc(S(=O)(=O)N[C@@H]2CCCC[C@H]2CN3C=NC4=CC=CC=C4N=C3N)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	446.8 (M + H)	4.29

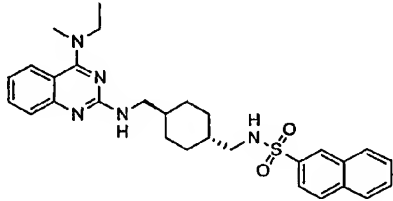
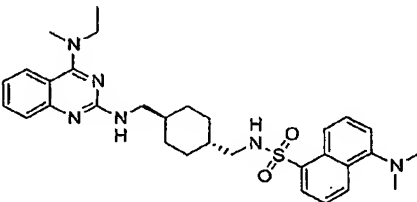
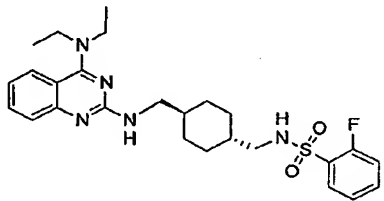
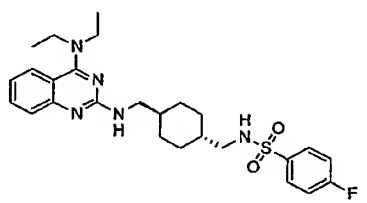
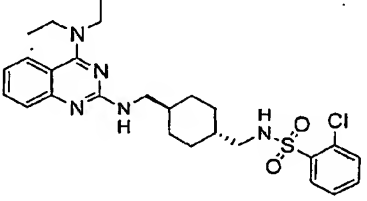
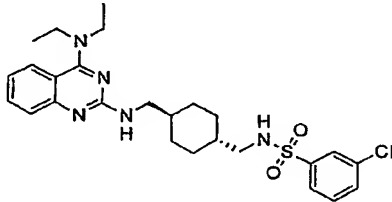
Example No.	Structure	ESI-MS	Retention Time (min)
2597	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4ccc(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.8 (M + H)	4.47
2598	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4cccc(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.8 (M + H)	4.53
2599	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4cccc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.6 (M + H)	4.55
2600	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4ccc(Cl)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	487.6 (M + H)	4.65
2601	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4cc(C(F)(F)F)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	556.6 (M + H)	4.91
2602	 <chem>CC1=NC2=CC=CC=C2N1NCC3CCCCC3NS(=O)(=O)c4ccccc4Br</chem> $\text{CF}_3\text{CO}_2\text{H}$	532.4 (M + H)	4.61

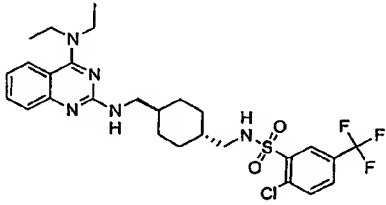
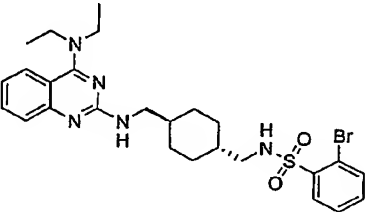
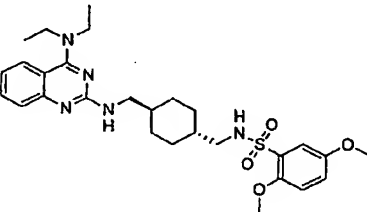
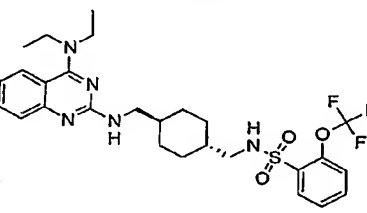
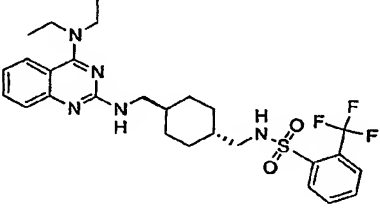
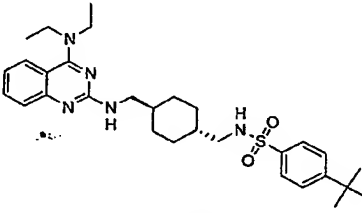
Example No.	Structure	ESI-MS	Retention Time (min)
2603	 <chem>CC1=NC2=C(N1)N=CN=C2CNC3CCCCC3NS(=O)(=O)c4cc(OC)cc(OC)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	514.8 (M + H)	4.43
2604	 <chem>CC1=NC2=C(N1)N=CN=C2CNC3CCCCC3NS(=O)(=O)c4cc(C(F)(F)F)ccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	538.6 (M + H)	4.80
2605	 <chem>CC1=NC2=C(N1)N=CN=C2CNC3CCCCC3NS(=O)(=O)c4ccc(C(C)(C)C)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	510.6 (M + H)	5.00
2606	 <chem>CC1=NC2=C(N1)N=CN=C2CNC3CCCCC3NS(=O)(=O)c4ccsc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	460.6 (M + H)	4.40
2607	 <chem>CCN(CC)C1=NC2=CC=CC=C2N=CN1CNC3CCCCC3NS(=O)(=O)c4cc(F)ccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.6 (M + H)	4.60
2608	 <chem>CCN(CC)C1=NC2=CC=CC=C2N=CN1CNC3CCCCC3NS(=O)(=O)c4ccc(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	484.6 (M + H)	4.64

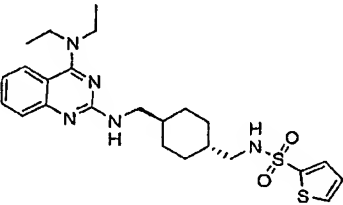
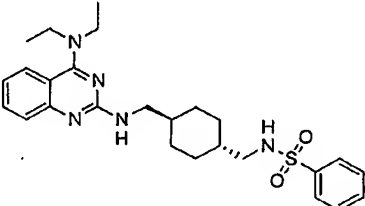
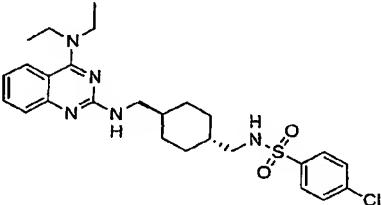
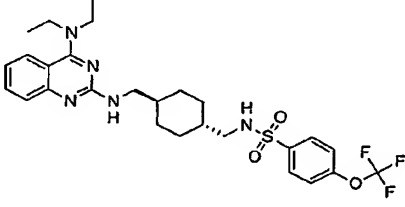
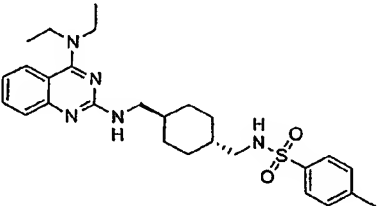
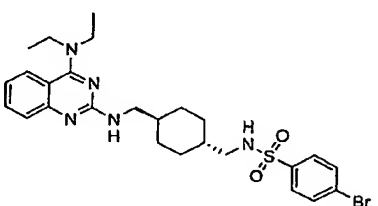
Example No.	Structure	ESI-MS	Retention Time (min)
2609	 $\text{CF}_3\text{CO}_2\text{H}$	503.6 (M + H)	4.74
2610	 $\text{CF}_3\text{CO}_2\text{H}$	502.6 (M + H)	4.86
2611	 $\text{CF}_3\text{CO}_2\text{H}$	570.8 (M + H)	5.00
2612	 $\text{CF}_3\text{CO}_2\text{H}$	546.0 (M + H)	4.80
2613	 $\text{CF}_3\text{CO}_2\text{H}$	528.8 (M + H)	4.63
2614	 $\text{CF}_3\text{CO}_2\text{H}$	552.8 (M + H)	4.90

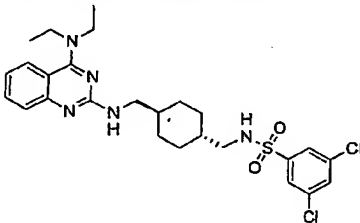
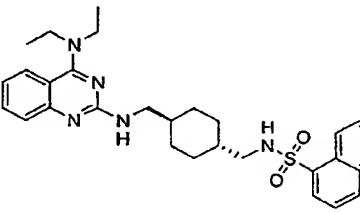
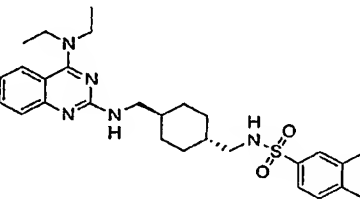
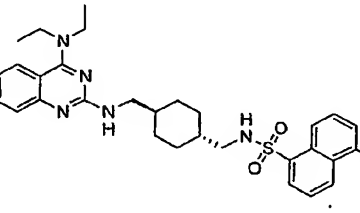
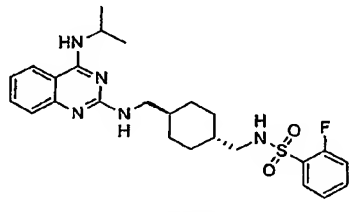
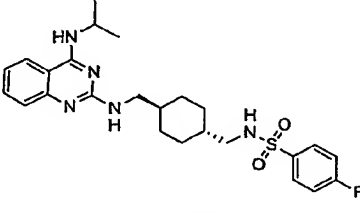
Example No.	Structure	ESI-MS	Retention Time (min)
2615	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)CC(F)(F)F)c3ccccc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	536.6 (M + H)	4.82
2616	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)Cc3ccc(C(C)(C)C)cc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	524.8 (M + H)	5.07
2617	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccsc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.6 (M + H)	4.55
2618	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccccc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	4.59
2619	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(Cl)cc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	502.6 (M + H)	4.81
2620	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(OC(F)(F)F)cc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	552.8 (M + H)	4.94

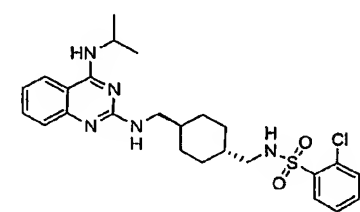
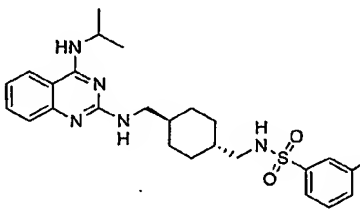
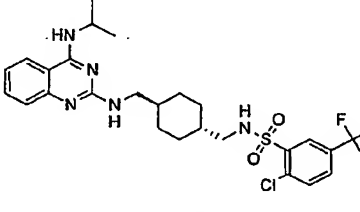
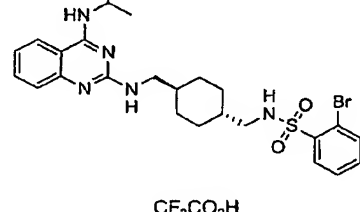
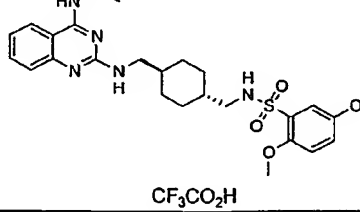
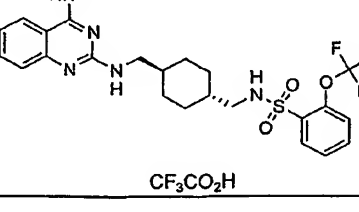
Example No.	Structure	ESI-MS	Retention Time (min)
2621	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccc(C)cc4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	482.6 (M + H)	4.73
2622	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccc(Br)cc4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	546.6 (M + H)	4.85
2623	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(Cl)cc(Cl)c4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	5.08
2624	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(OC(F)(F)F)cc(Br)c4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	630.4 (M + H)	5.11
2625	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(C(F)(F)F)cc(C(F)(F)F)c4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	604.6 (M + H)	5.16
2626	 <chem>CCN(C)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4c5ccccc5cc4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.75

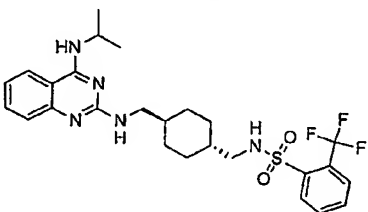
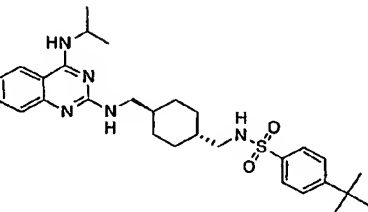
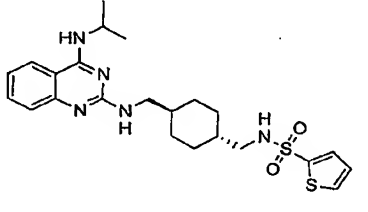
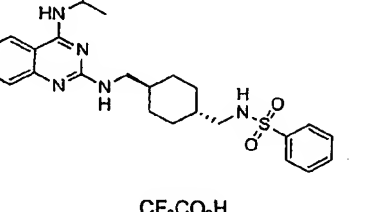
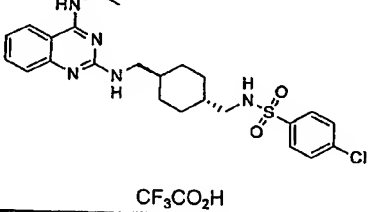
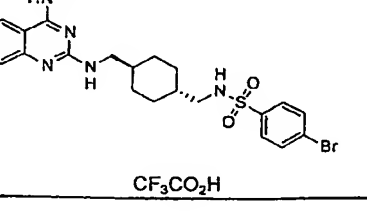
Example No.	Structure	ESI-MS	Retention Time (min)
2627	 $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.91
2628	 $2\text{CF}_3\text{CO}_2\text{H}$	561.6 (M + H)	4.61
2629	 $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.75
2630	 $\text{CF}_3\text{CO}_2\text{H}$	500.2 (M + H)	4.85
2631	 $\text{CF}_3\text{CO}_2\text{H}$	516.6 (M + H)	4.81
2632	 $\text{CF}_3\text{CO}_2\text{H}$	516.6 (M + H)	4.95

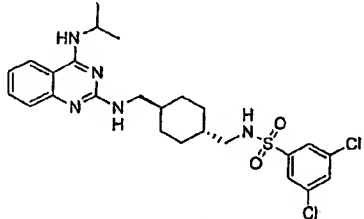
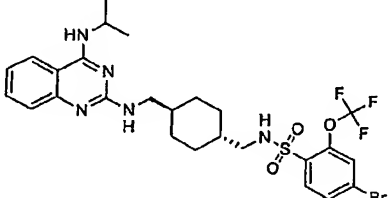
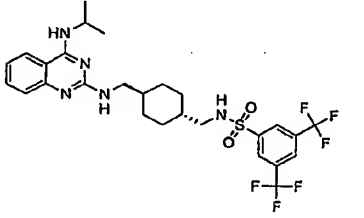
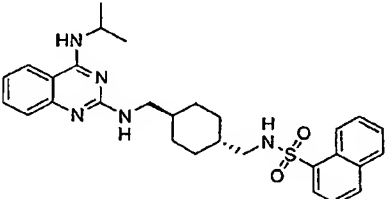
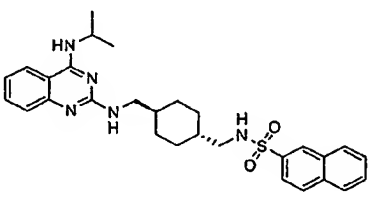
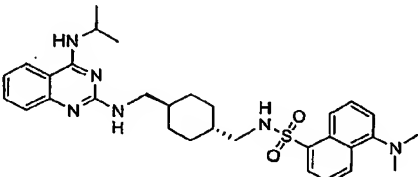
Example No.	Structure	ESI-MS	Retention Time (min)
2633	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(F)(F)Fcc4Cl)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	584.6 (M + H)	5.18
2634	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccccc4Br)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	560.6 (M + H)	4.87
2635	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(OC)cc4OC)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	542.8 (M + H)	4.80
2636	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccccc4OC(F)(F)F)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	566.6 (M + H)	5.01
2637	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccccc4C(F)(F)F)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.8 (M + H)	4.95
2638	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccc(C(C)(C)C)cc4)C3CCCCC3</chem> $\text{CF}_3\text{CO}_2\text{H}$	538.6 (M + H)	5.20

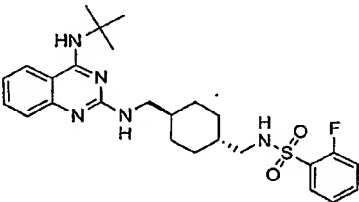
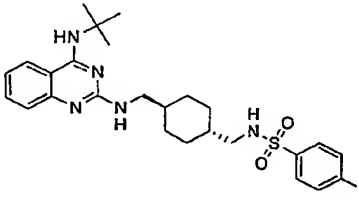
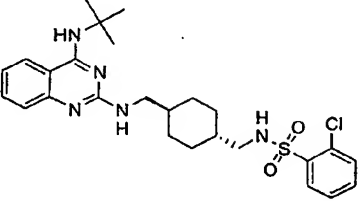
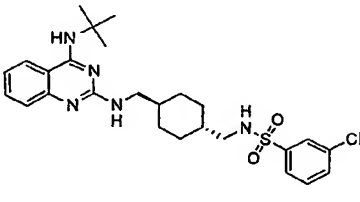
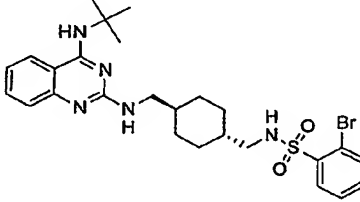
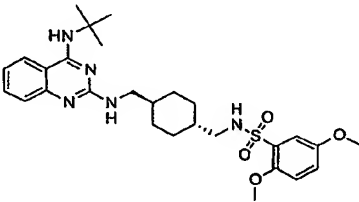
Example No.	Structure	ESI-MS	Retention Time (min)
2639	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccsc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.6 (M + H)	4.65
2640	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccccc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	482.6 (M + H)	4.73
2641	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(Cl)cc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.8 (M + H)	4.97
2642	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(OC(F)(F)F)cc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	566.6 (M + H)	5.12
2643	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(C)cc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	496.8 (M + H)	4.89
2644	 <chem>CCN(CC)c1nc2c(ncn2C1CCN(S(=O)(=O)c3ccc(Br)cc3)C4CCCCC4)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	560.0 (M + H)	4.98

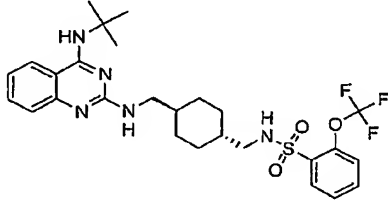
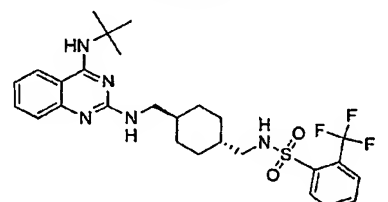
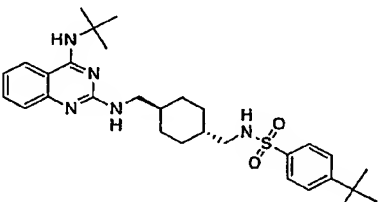
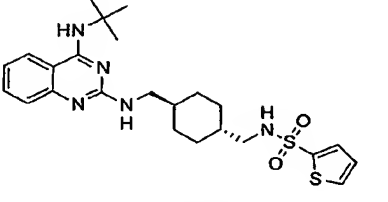
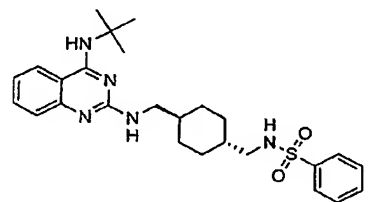
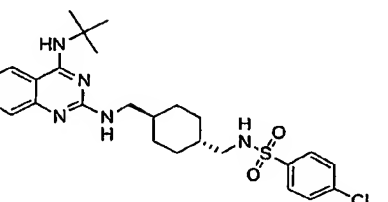
Example No.	Structure	ESI-MS	Retention Time (min)
2645	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4cc(Cl)cc(Cl)c4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.6 (M + H)	5.21
2646	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4c5ccccc5c6ccccc46)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	532.6 (M + H)	4.99
2647	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4c5ccccc5c6ccccc46)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	532.6 (M + H)	5.03
2648	 <chem>CCN(CC)c1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4c5ccccc5c6ccccc46N(C)C)C3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	575.8 (M + H)	4.80
2649	 <chem>CC(C)Nc1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccccc4F)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.6 (M + H)	4.64
2650	 <chem>CC(C)Nc1nc2c(ncn2C1CC3CCCCC3NS(=O)(=O)c4ccc(F)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.6 (M + H)	4.66

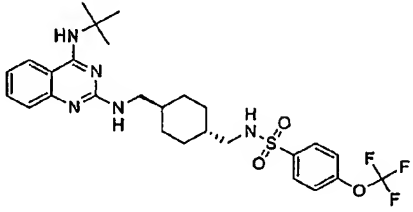
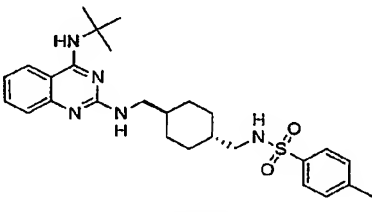
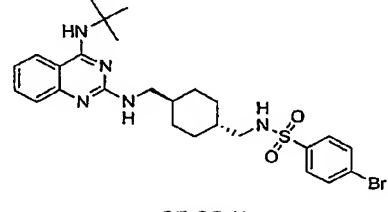
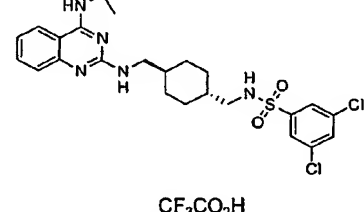
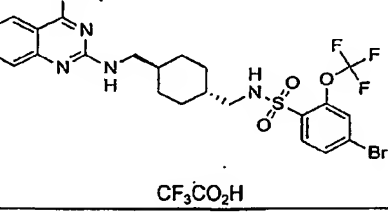
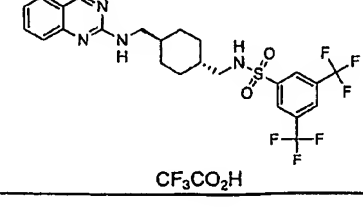
Example No.	Structure	ESI-MS	Retention Time (min)
2651	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3ccccc3Cl)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	502.6 (M + H)	4.72
2652	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3cccc(c3)Cl)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	502.6 (M + H)	4.87
2653	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3cc(C(F)(F)F)ccc3Cl)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	570.6 (M + H)	5.03
2654	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3ccccc3Br)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	546.6 (M + H)	4.77
2655	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3cc(OC)cc(OC)c3)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	528.8 (M + H)	4.68
2656	 <chem>CC(C)Nc1nc2c(ncn2C1CCN(C1)CS(=O)(=O)c3cc(OC(F)(F)F)ccc3)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	552.8 (M + H)	4.89

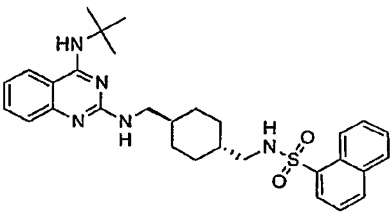
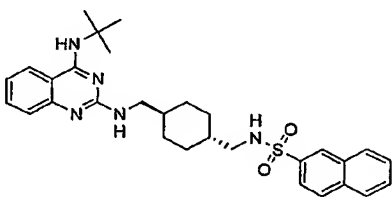
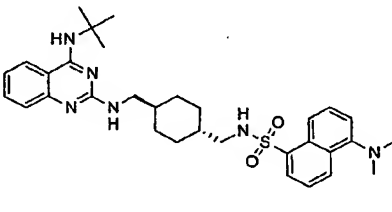
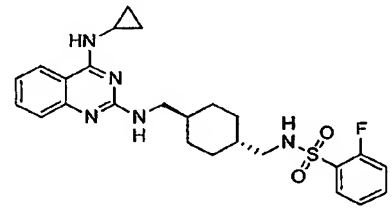
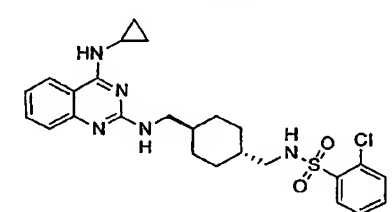
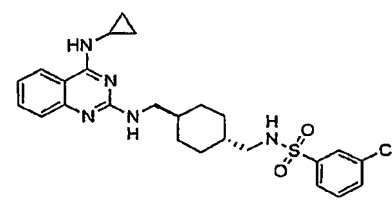
Example No.	Structure	ESI-MS	Retention Time (min)
2657	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3cc(F)(F)Fcc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	536.6 (M + H)	4.85
2658	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3ccc(C(C)(C)C)cc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	524.8 (M + H)	5.15
2659	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3ccsc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.8 (M + H)	4.63
2660	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3ccccc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	4.61
2661	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3ccc(Cl)cc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	502.6 (M + H)	4.86
2662	 <chem>CC(C)Nc1nc2c(ncn2C1CCCCC1NS(=O)(=O)c3ccc(Br)cc3)cc3ccccc33</chem> $\text{CF}_3\text{CO}_2\text{H}$	546.6 (M + H)	4.64

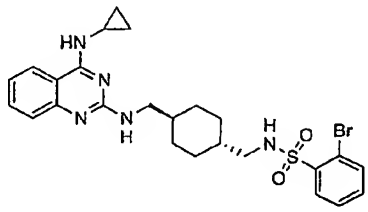
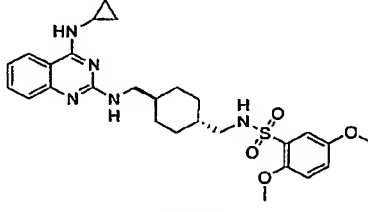
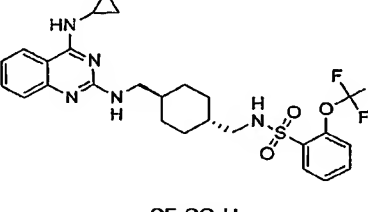
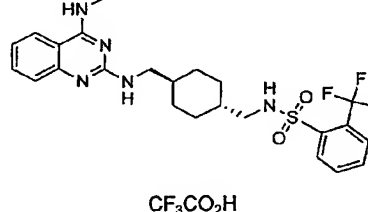
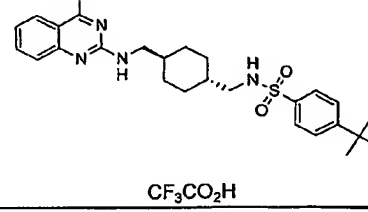
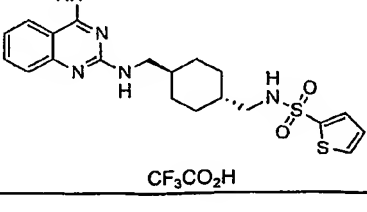
Example No.	Structure	ESI-MS	Retention Time (min)
2663	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc(Cl)cc(Cl)c3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	4.81
2664	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc(Br)cc(OC(F)(F)F)c3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	630.4 (M + H)	4.85
2665	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc(C(F)(F)F)cc(C(F)(F)F)c3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	604.6 (M + H)	4.87
2666	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc4ccccc4cc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.67
2667	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc4ccccc4cc3)cc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.90
2668	 <chem>CC(C)Nc1nc2c(ncn2C1CCN1S(=O)(=O)c3cc4ccccc4c(N(C)C)c3)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	561.6 (M + H)	4.64

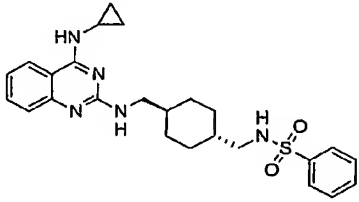
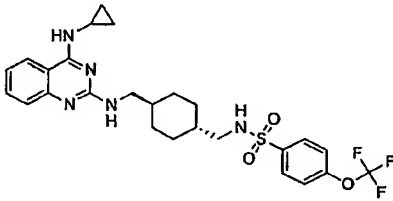
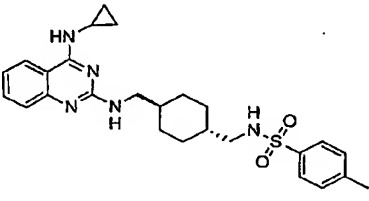
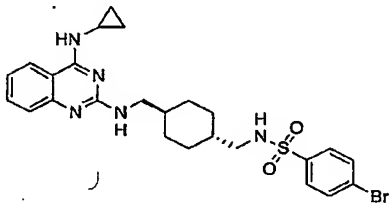
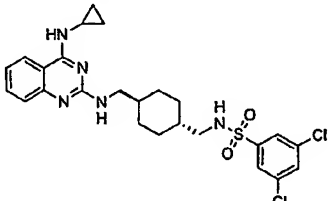
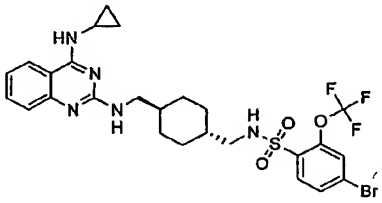
Example No.	Structure	ESI-MS	Retention Time (min)
2669	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4ccccc4F</chem> $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.73
2670	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4ccc(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.74
2671	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4ccccc4Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.6 (M + H)	4.89
2672	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4cccc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.6 (M + H)	4.93
2673	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4ccccc4Br</chem> $\text{CF}_3\text{CO}_2\text{H}$	560.0 (M + H)	4.89
2674	 <chem>CC1(C)C(C(C1)N2C=NC3C=CC=CC=C3N2C)S(=O)(=O)c4cc(OC)c(OC)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	542.8 (M + H)	4.76

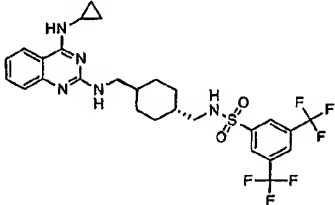
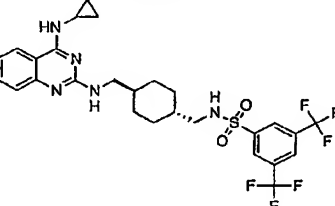
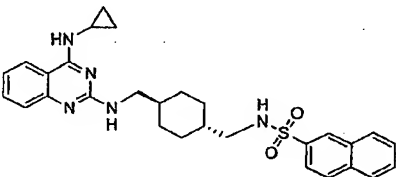
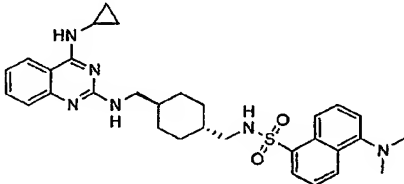
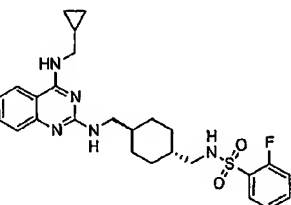
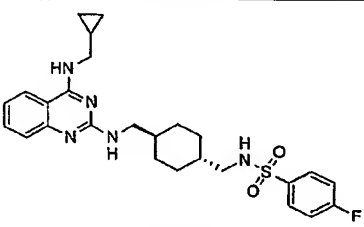
Example No.	Structure	ESI-MS	Retention Time (min)
2675	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)OC(F)(F)F)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	566.6 (M + H)	5.03
2676	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)C(F)(F)F)c3ccccc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.8 (M + H)	4.96
2677	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccc(C(C)(C)C)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	538.8 (M + H)	5.25
2678	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccsc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.6 (M + H)	4.67
2679	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	482.4 (M + H)	4.71
2680	 <chem>CC(C)(C)Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccc(Cl)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.6 (M + H)	4.95

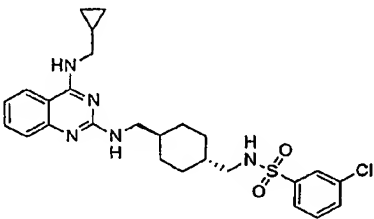
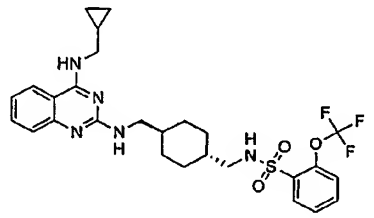
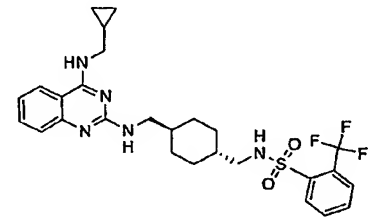
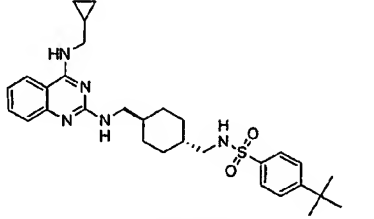
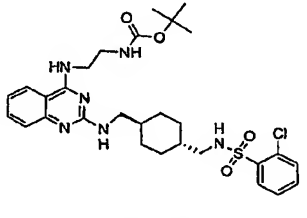
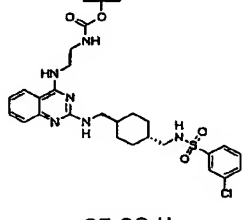
Example No.	Structure	ESI-MS	Retention Time (min)
2681	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1ccc(OC(F)(F)F)cc1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	566.8 (M + H)	5.07
2682	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1ccc(C)cc1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	496.8 (M + H)	4.83
2683	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1ccc(Br)cc1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	560.6 (M + H)	5.01
2684	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1cc(Cl)cc(Cl)c1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.6 (M + H)	5.07
2685	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1cc(Br)cc(OC(F)(F)F)c1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	644.6 (M + H)	5.29
2686	 <chem>CC1(C)NC2=NC(=NC3=CC=CC=C3N2)NC(C1)CCN(S(=O)(=O)c1cc(C(F)(F)F)cc(C(F)(F)F)c1)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	618.6 (M + H)	5.25

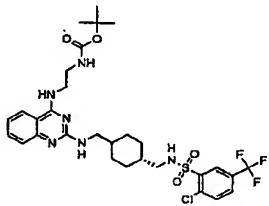
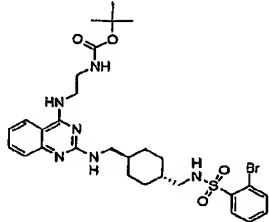
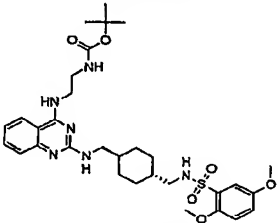
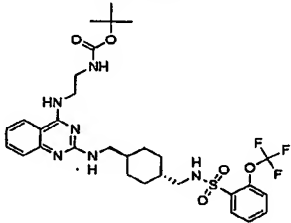
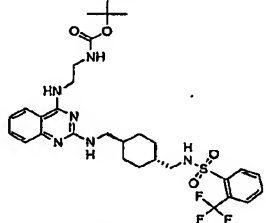
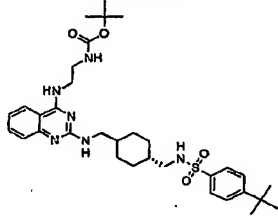
Example No.	Structure	ESI-MS	Retention Time (min)
2687	 $\text{CF}_3\text{CO}_2\text{H}$	532.6 (M + H)	5.01
2688	 $\text{CF}_3\text{CO}_2\text{H}$	532.6 (M + H)	5.04
2689	 $2\text{CF}_3\text{CO}_2\text{H}$	575.8 (M + H)	4.75
2690	 $\text{CF}_3\text{CO}_2\text{H}$	484.6 (M + H)	4.51
2691	 $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.59
2692	 $\text{CF}_3\text{CO}_2\text{H}$	500.8 (M + H)	4.71

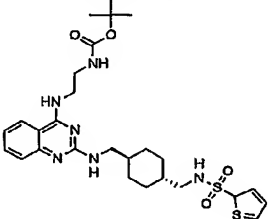
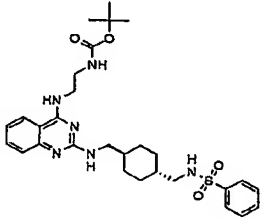
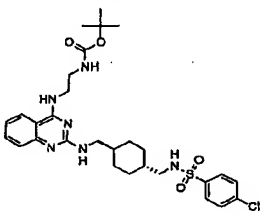
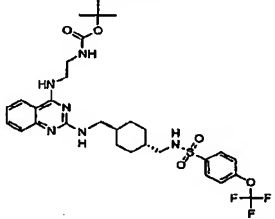
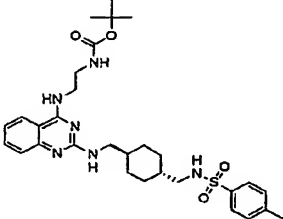
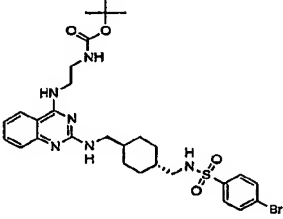
Example No.	Structure	ESI-MS	Retention Time (min)
2693	 <chem>BrC1=CC=C(C=C1)S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)C5CC5</chem> $\text{CF}_3\text{CO}_2\text{H}$	544.6 (M + H)	4.63
2694	 <chem>COc1cc(OC)c(S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	526.8 (M + H)	4.55
2695	 <chem>FC(F)(F)c1ccc(S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.6 (M + H)	4.79
2696	 <chem>FC(F)(F)c1ccc(S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	534.6 (M + H)	4.69
2697	 <chem>CC(C)(C)c1ccc(S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	522.4 (M + H)	5.03
2698	 <chem>c1cc(S(=O)(=O)N[C@H]2CCCC[C@H]2CNc3nc4c(ncn4C3)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.8 (M + H)	4.43

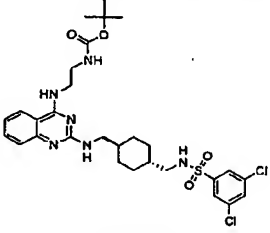
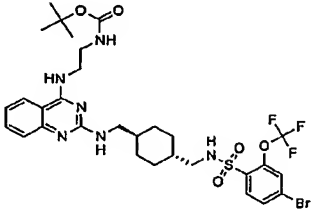
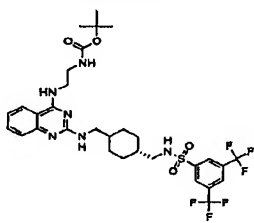
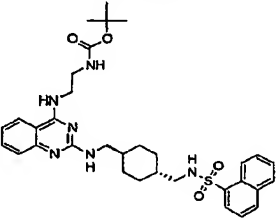
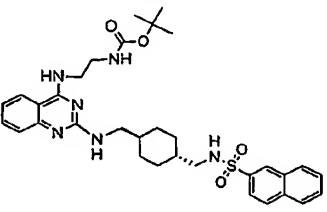
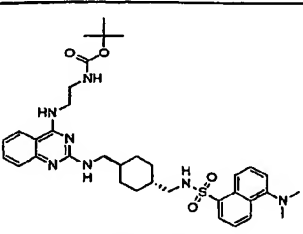
Example No.	Structure	ESI-MS	Retention Time (min)
2699	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1ccccc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	466.6 (M + H)	4.50
2700	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1ccc(OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	550.6 (M + H)	4.87
2701	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1ccc(C)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.6 (M + H)	4.65
2702	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1ccc(Br)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	544.6 (M + H)	4.75
2703	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1cc(Cl)cc(Cl)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	534.6 (M + H)	4.90
2704	 <chem>CC1(C)Nc2nc3c(ncn3C2CC1)cccc2S(=O)(=O)c1cc(Br)cc(OC(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	628.6 (M + H)	5.08

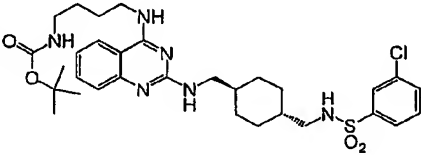
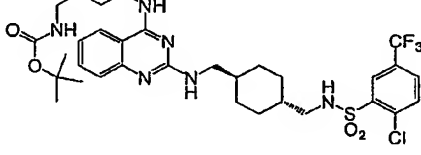
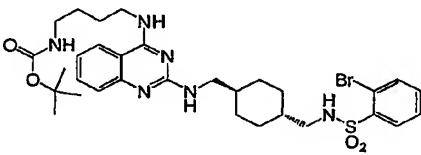
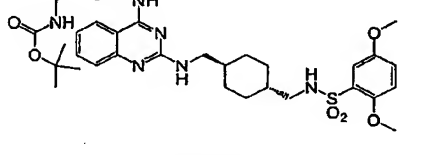
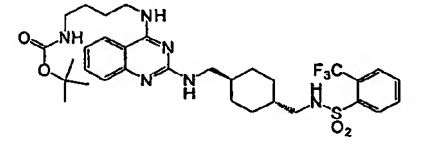
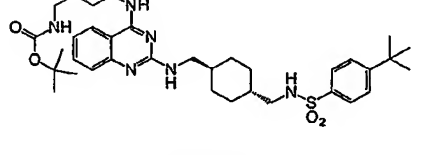
Example No.	Structure	ESI-MS	Retention Time (min)
2705	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc(C(F)(F)F)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	602.6 (M + H)	5.10
2706	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc(C(F)(F)F)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.8 (M + H)	4.71
2707	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc3ccccc3c2</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.8 (M + H)	4.81
2708	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc3c(c2)N(C)ccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	559.6 (M + H)	4.50
2709	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc(F)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	498.8 (M + H)	4.64
2710	 <chem>CC1(C)CCC(CC1)NS(=O)(=O)c2ccc(F)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	498.8 (M + H)	4.73

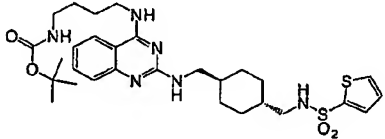
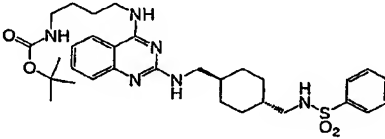
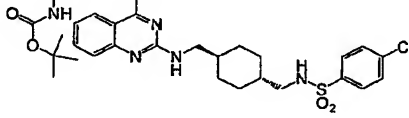
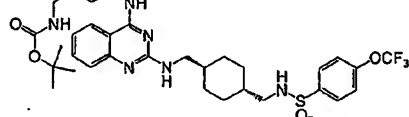
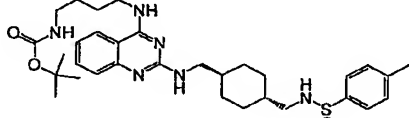
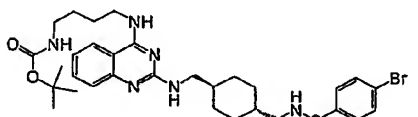
Example No.	Structure	ESI-MS	Retention Time (min)
2711	 $\text{CF}_3\text{CO}_2\text{H}$	514.8 (M + H)	4.87
2712	 $\text{CF}_3\text{CO}_2\text{H}$	564.6 (M + H)	4.93
2713	 $\text{CF}_3\text{CO}_2\text{H}$	548.6 (M + H)	4.87
2714	 $\text{CF}_3\text{CO}_2\text{H}$	536.6 (M + H)	5.19
2715	 $\text{CF}_3\text{CO}_2\text{H}$	603.8 (M + H)	4.76
2716	 $\text{CF}_3\text{CO}_2\text{H}$	603.4 (M + H)	4.87

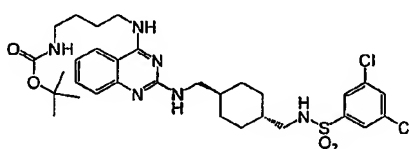
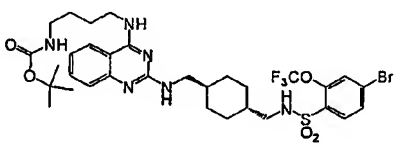
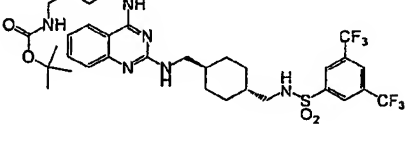
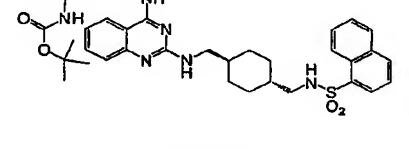
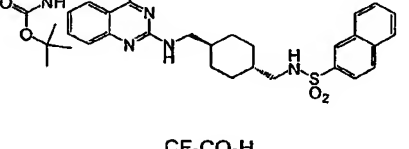
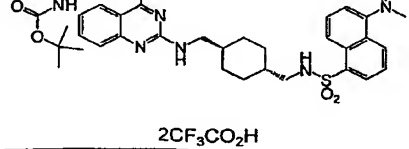
Example No.	Structure	ESI-MS	Retention Time (min)
2717	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4cc(Cl)cc(C(F)(F)F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	671.6 (M + H)	5.05
2718	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4cc(Br)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	647.6 (M + H)	4.79
2719	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4cc(OC)c(OC)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	629.8 (M + H)	4.67
2720	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4cc(OC(F)(F)F)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	653.8 (M + H)	4.91
2721	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4cc(C(F)(F)F)c(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	637.8 (M + H)	4.85
2722	 <chem>CC(C)(C)OC(=O)NCCNc1nc2ccccc2n1CNC3CCCCC3NS(=O)(=O)c4ccc(C(C)(C)C)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	625.8 (M + H)	5.14

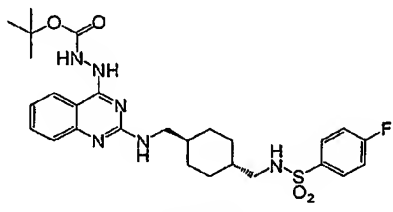
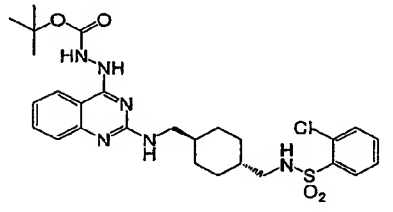
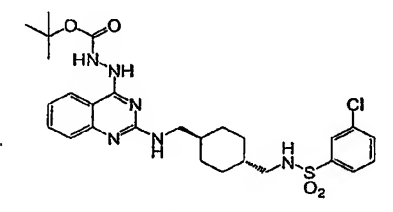
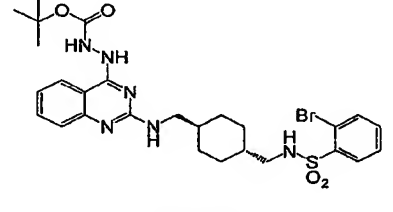
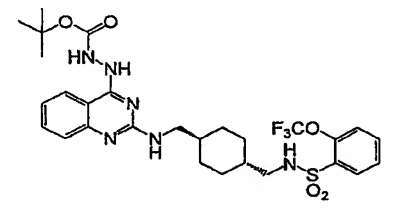
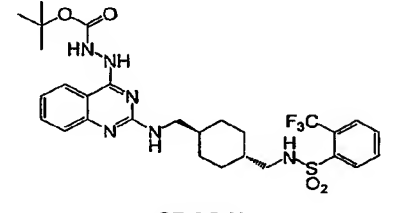
Example No.	Structure	ESI-MS	Retention Time (min)
2723	 $\text{CF}_3\text{CO}_2\text{H}$	575.6 (M + H)	4.63
2724	 $\text{CF}_3\text{CO}_2\text{H}$	569.8 (M + H)	4.66
2725	 $\text{CF}_3\text{CO}_2\text{H}$	603.8 (M + H)	4.88
2726	 $\text{CF}_3\text{CO}_2\text{H}$	653.8 (M + H)	5.01
2727	 $\text{CF}_3\text{CO}_2\text{H}$	583.8 (M + H)	4.77
2728	 $\text{CF}_3\text{CO}_2\text{H}$	647 (M + H)	4.92

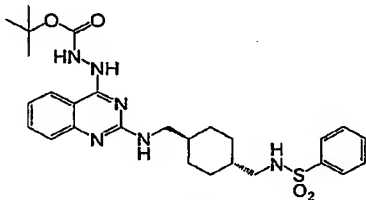
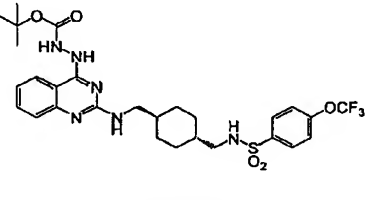
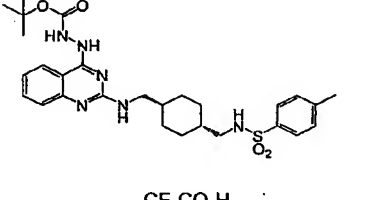
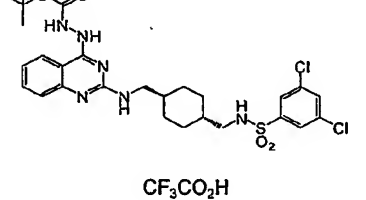
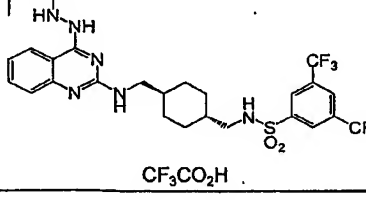
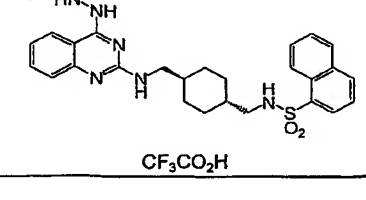
Example No.	Structure	ESI-MS	Retention Time (min)
2729	 $\text{CF}_3\text{CO}_2\text{H}$	637.8 (M + H)	5.13
2730	 $\text{CF}_3\text{CO}_2\text{H}$	731.6 (M + H)	5.19
2731	 $\text{CF}_3\text{CO}_2\text{H}$	705.8 (M + H)	5.22
2732	 $\text{CF}_3\text{CO}_2\text{H}$	619.8 (M + H)	4.91
2733	 $\text{CF}_3\text{CO}_2\text{H}$	619.8 (M + H)	4.93
2734	 $2\text{CF}_3\text{CO}_2\text{H}$	663.0 (M + H)	4.67

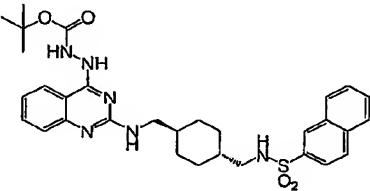
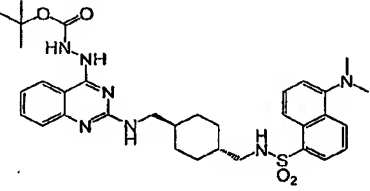
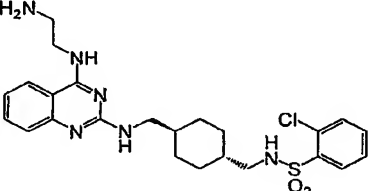
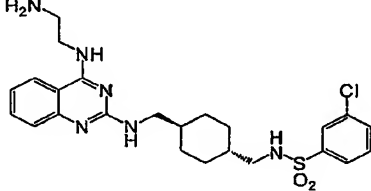
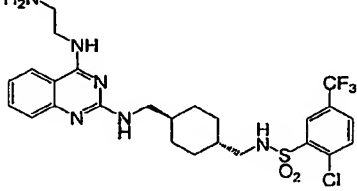
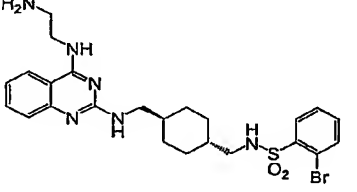
Example No.	Structure	ESI-MS	Retention Time (min)
2735	 $\text{CF}_3\text{CO}_2\text{H}$	631.8 (M + H)	5.01
2736	 $\text{CF}_3\text{CO}_2\text{H}$	699.0 (M + H)	5.19
2737	 $\text{CF}_3\text{CO}_2\text{H}$	675.8 (M + H)	4.95
2738	 $\text{CF}_3\text{CO}_2\text{H}$	657.8 (M + H)	4.81
2739	 $\text{CF}_3\text{CO}_2\text{H}$	665.8 (M + H)	4.97
2740	 $\text{CF}_3\text{CO}_2\text{H}$	653.8 (M + H)	5.27

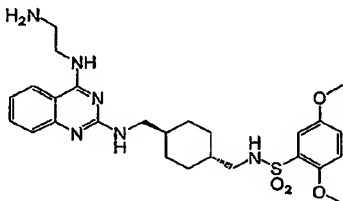
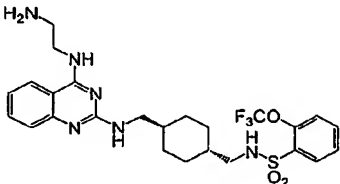
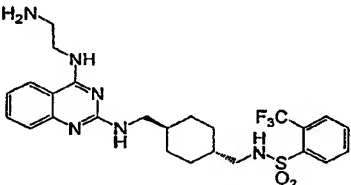
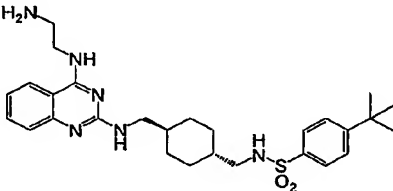
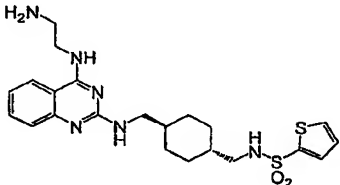
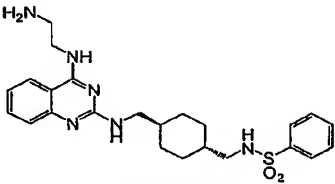
Example No.	Structure	ESI-MS	Retention Time (min)
2741	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccsc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	603.4 (M + H)	4.77
2742	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	597.8 (M + H)	4.79
2743	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccc(Cl)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	631.8 (M + H)	5.02
2744	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccc(OC(F)(F)F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	681.8 (M + H)	5.14
2745	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccc(C)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	611.8 (M + H)	4.93
2746	 <chem>CC(C)(C)OC(=O)N[C@@H]1C=CC2=C(N1)N=CN=C2N[C@@H]3CCCC[C@H]3CNS(=O)(=O)c4ccc(Br)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	675.0 (M + H)	5.05

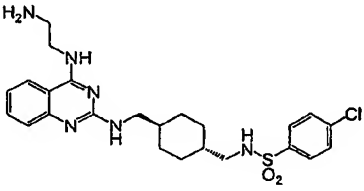
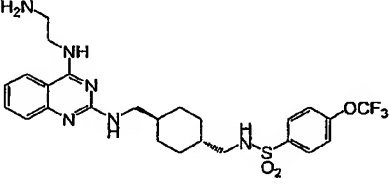
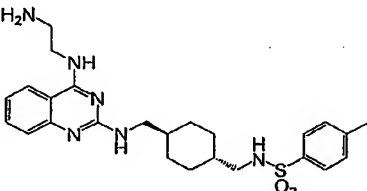
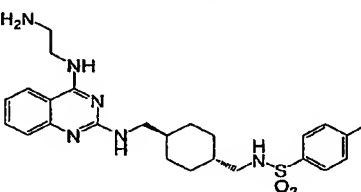
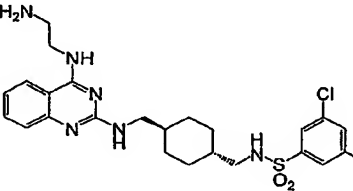
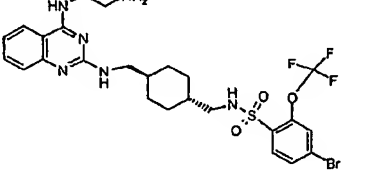
Example No.	Structure	ESI-MS	Retention Time (min)
2747	 $\text{CF}_3\text{CO}_2\text{H}$	665.8 (M + H)	5.29
2748	 $\text{CF}_3\text{CO}_2\text{H}$	759.6 (M + H)	5.31
2749	 $\text{CF}_3\text{CO}_2\text{H}$	733.8 (M + H)	5.36
2750	 $\text{CF}_3\text{CO}_2\text{H}$	647.8 (M + H)	5.05
2751	 $\text{CF}_3\text{CO}_2\text{H}$	647.8 (M + H)	5.08
2752	 $2\text{CF}_3\text{CO}_2\text{H}$	691.0 (M + H)	4.89

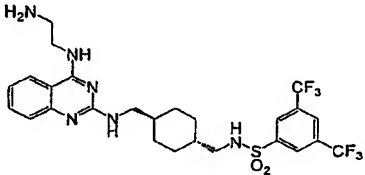
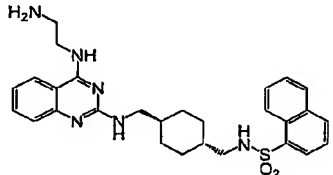
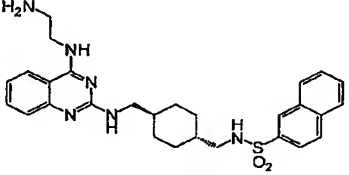
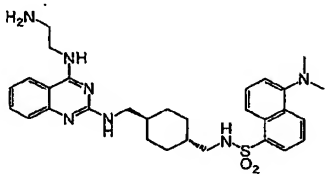
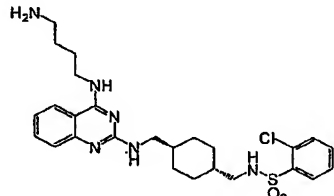
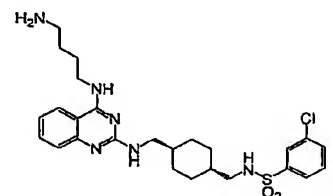
Example No.	Structure	ESI-MS	Retention Time (min)
2753	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	559.6 (M + H)	4.51
2754	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(Cl)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	575.6 (M + H)	4.57
2755	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(Cl)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	575.6 (M + H)	4.69
2756	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(Br)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	619.6 (M + H)	4.63
2757	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(OC(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	625.8 (M + H)	4.72
2758	 <chem>CC(C)(C)OC(=O)NNc1nc2c(nc1)nc3ccccc23CNC4CCCCC4CNS(=O)(=O)c5ccc(C(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	609.8 (M + H)	4.67

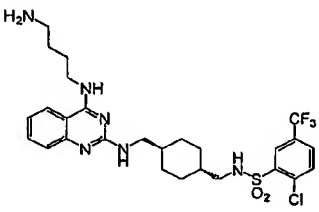
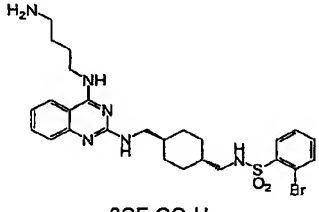
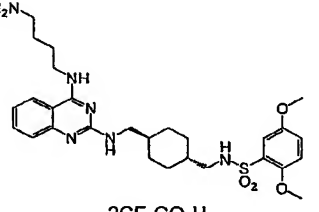
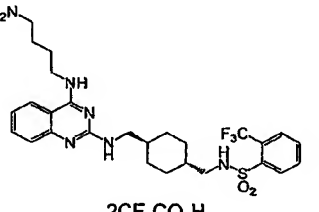
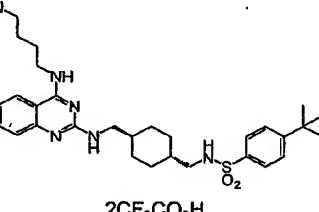
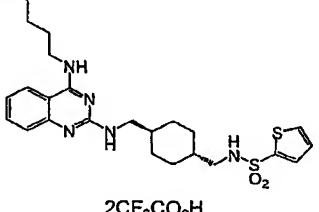
Example No.	Structure	ESI-MS	Retention Time (min)
2759	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3ccccc3)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	541.8 (M + H)	4.45
2760	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3ccc(OC(F)(F)F)cc3)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	625.8 (M + H)	4.38
2761	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3ccc(C)cc3)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	555.8 (M + H)	4.57
2762	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3cc(Cl)cc(Cl)c3)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	609.8 (M + H)	4.94
2763	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3cc(C(F)(F)F)cc(C(F)(F)F)c3)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	677.8 (M + H)	5.05
2764	 <chem>CC(C)(C)OC(=O)NNc1nc2c(ncn2C1CCN(C1)CNS(=O)(=O)c3cccc4ccccc34)nc3ccccc13</chem> $\text{CF}_3\text{CO}_2\text{H}$	591.6 (M + H)	4.73

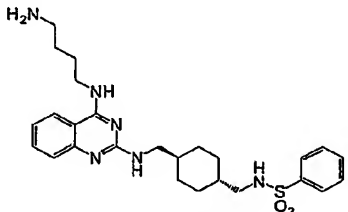
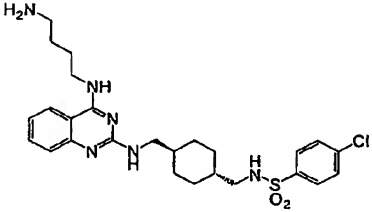
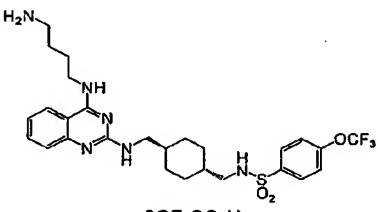
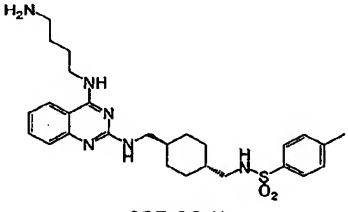
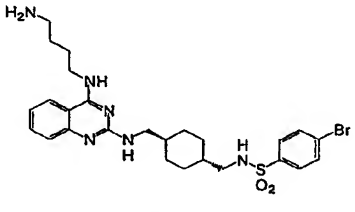
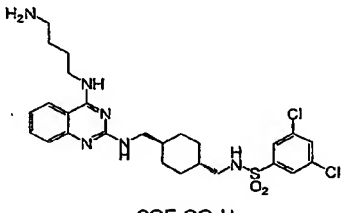
Example No.	Structure	ESI-MS	Retention Time (min)
2765	 <chem>CC(C)(C)OC(=O)NNc1nc2ccccc2n1CNC3CCCCC3CNS(=O)(=O)c4ccc5ccccc45</chem> $\text{CF}_3\text{CO}_2\text{H}$	591.6 (M + H)	4.75
2766	 <chem>CC(C)(C)OC(=O)NNc1nc2ccccc2n1CNC3CCCCC3CNS(=O)(=O)c4ccc5c(c4)CN(C)CC5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	635.0 (M + H)	4.47
2767	 <chem>NCCNC1=NC2=CC=CC=C2N=C1CNC3CCCCC3CNS(=O)(=O)c4ccccc4Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	503.6 (M + H)	3.83
2768	 <chem>NCCNC1=NC2=CC=CC=C2N=C1CNC3CCCCC3CNS(=O)(=O)c4ccc(Cl)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	503.6 (M + H)	3.99
2769	 <chem>NCCNC1=NC2=CC=CC=C2N=C1CNC3CCCCC3CNS(=O)(=O)c4cc(Cl)ccc4C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	571.6 (M + H)	4.16
2770	 <chem>NCCNC1=NC2=CC=CC=C2N=C1CNC3CCCCC3CNS(=O)(=O)c4ccccc4Br</chem> $2\text{CF}_3\text{CO}_2\text{H}$	547.6 (M + H)	3.85

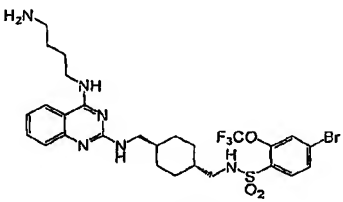
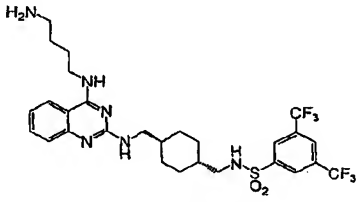
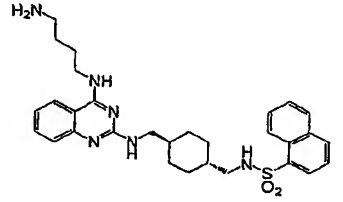
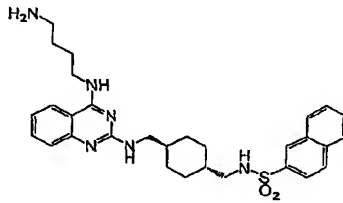
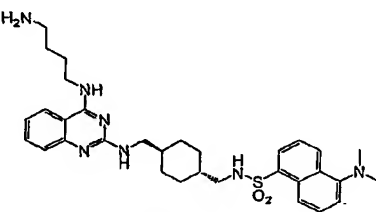
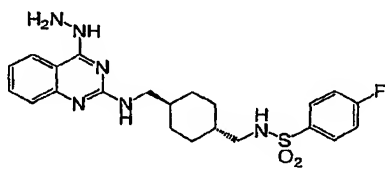
Example No.	Structure	ESI-MS	Retention Time (min)
2771	 <chem>2CF3CO2H</chem>	529.6 (M + H)	3.75
2772	 <chem>2CF3CO2H</chem>	553.8 (M + H)	3.99
2773	 <chem>2CF3CO2H</chem>	537.6 (M + H)	3.93
2774	 <chem>2CF3CO2H</chem>	525.8 (M + H)	4.22
2775	 <chem>2CF3CO2H</chem>	475.6 (M + H)	3.64
2776	 <chem>2CF3CO2H</chem>	469.6 (M + H)	3.71

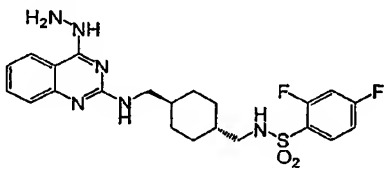
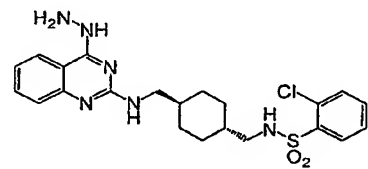
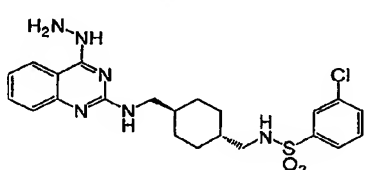
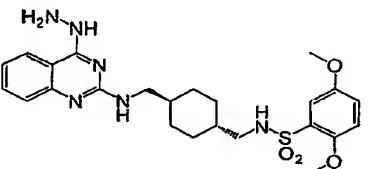
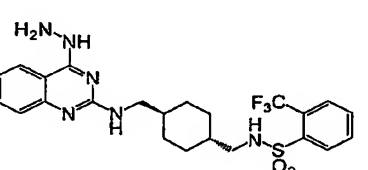
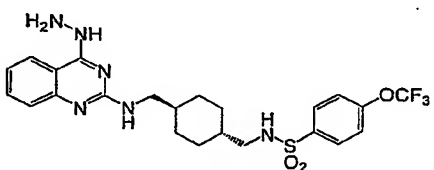
Example No.	Structure	ESI-MS	Retention Time (min)
2777	 2CF ₃ CO ₂ H	503.6 (M + H)	3.97
2778	 2CF ₃ CO ₂ H	553.8 (M + H)	4.17
2779	 2CF ₃ CO ₂ H	483.4 (M + H)	3.87
2780	 2CF ₃ CO ₂ H	547.6 (M + H)	4.04
2781	 2CF ₃ CO ₂ H	537.4 (M + H)	4.23
2782	 2CF ₃ CO ₂ H	631.6 (M + H)	4.23

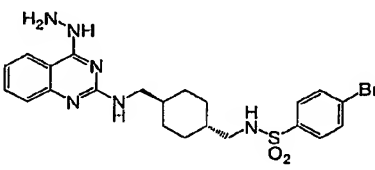
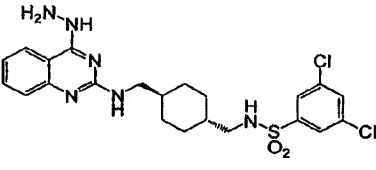
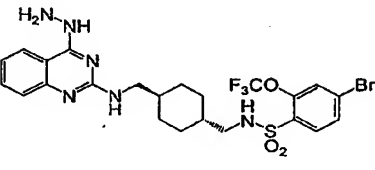
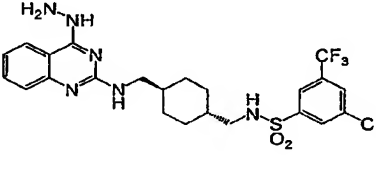
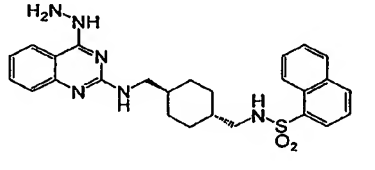
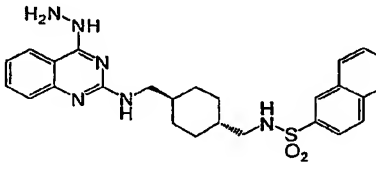
Example No.	Structure	ESI-MS	Retention Time (min)
2783	 <p>2CF₃CO₂H</p>	605.8 (M + H)	4.41
2784	 <p>2CF₃CO₂H</p>	519.6 (M + H)	4.01
2785	 <p>2CF₃CO₂H</p>	519.6 (M + H)	4.07
2786	 <p>3CF₃CO₂H</p>	562.6 (M + H)	3.77
2787	 <p>2CF₃CO₂H</p>	531.6 (M + H)	3.90
2788	 <p>2CF₃CO₂H</p>	531.6 (M + H)	4.04

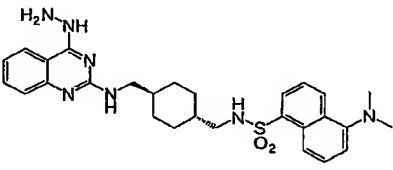
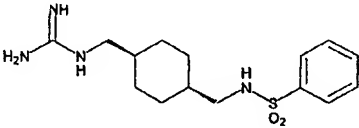
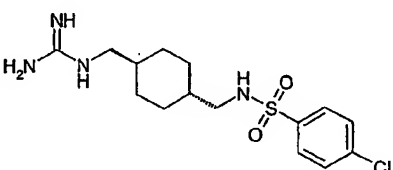
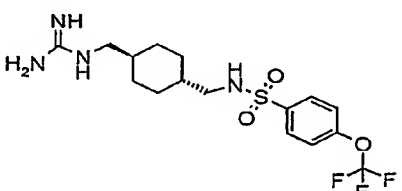
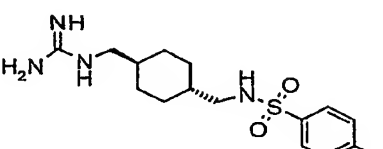
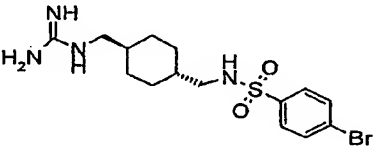
Example No.	Structure	ESI-MS	Retention Time (min)
2789	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4cc(Cl)cc(C(F)(F)F)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	599.6 (M + H)	4.24
2790	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4cc(Br)ccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	575.0 (M + H)	3.95
2791	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4cc(OC)cc(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	557.6 (M + H)	3.86
2792	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4cc(C(F)(F)F)ccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	565.6 (M + H)	4.03
2793	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4ccc(C(C)(C)C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	554 (M + H)	4.29
2794	 <chem>NCCCNc1nc2ccccc2n1CNC3CCCCC3CS(=O)(=O)c4ccsc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	503.6 (M + H)	3.78

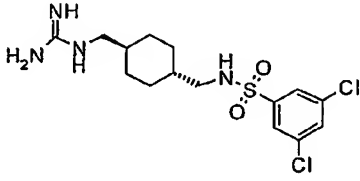
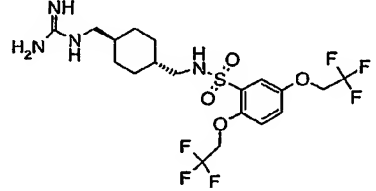
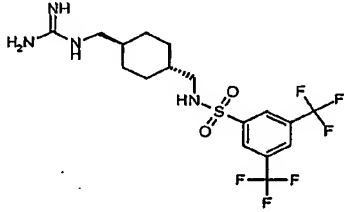
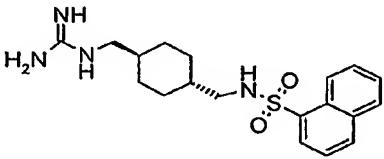
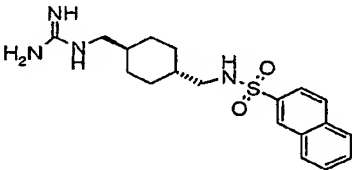
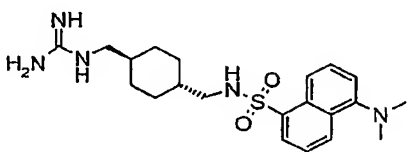
Example No.	Structure	ESI-MS	Retention Time (min)
2795	 $2\text{CF}_3\text{CO}_2\text{H}$	497.6 (M + H)	3.83
2796	 $2\text{CF}_3\text{CO}_2\text{H}$	531.6 (M + H)	4.05
2797	 $2\text{CF}_3\text{CO}_2\text{H}$	582.0 (M + H)	4.23
2798	 $2\text{CF}_3\text{CO}_2\text{H}$	511 (M + H)	3.95
2799	 $2\text{CF}_3\text{CO}_2\text{H}$	575.6 (M + H)	4.10
2800	 $2\text{CF}_3\text{CO}_2\text{H}$	565.0 (M + H)	4.32

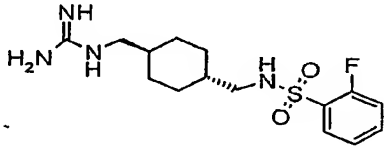
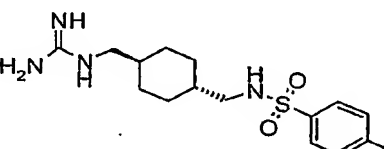
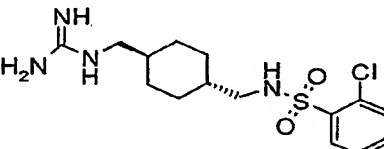
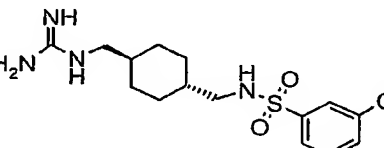
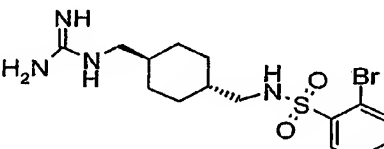
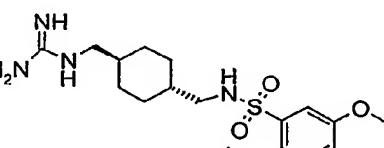
Example No.	Structure	ESI-MS	Retention Time (min)
2801	 $2\text{CF}_3\text{CO}_2\text{H}$	659.6 (M + H)	4.35
2802	 $2\text{CF}_3\text{CO}_2\text{H}$	634.0 (M + H)	4.43
2803	 $2\text{CF}_3\text{CO}_2\text{H}$	547.6 (M + H)	4.09
2804	 $2\text{CF}_3\text{CO}_2\text{H}$	547.6 (M + H)	4.15
2805	 $3\text{CF}_3\text{CO}_2\text{H}$	590.6 (M + H)	3.93
2806	 $2\text{CF}_3\text{CO}_2\text{H}$	459.6 (M + H)	4.07

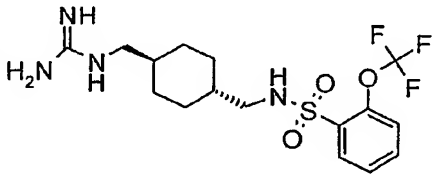
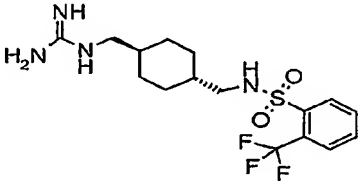
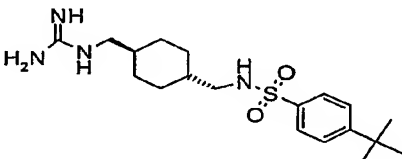
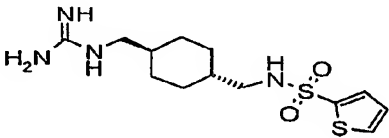
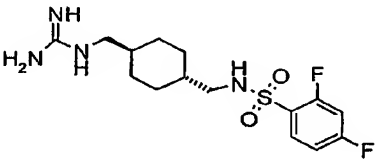
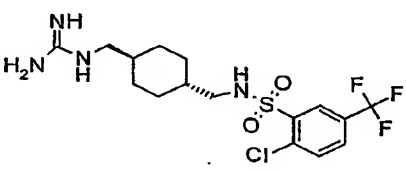
Example No.	Structure	ESI-MS	Retention Time (min)
2807	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4cc(F)cc(F)c4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	477.6 (M + H)	4.07
2808	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccccc4Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	475.6 (M + H)	4.07
2809	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4cccc(Cl)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	475.6 (M + H)	4.23
2810	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4cc(OC)cc(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	501.8 (M + H)	4.15
2811	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccccc4C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	509.4 (M + H)	4.27
2812	 <chem>Nc1nc2c(ncn2C1CCNCC3CCCCC3NS(=O)(=O)c4ccc(OC(F)(F)F)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	525.6 (M + H)	4.37

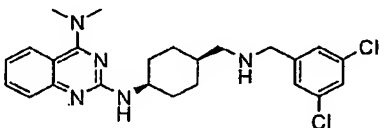
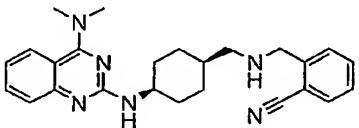
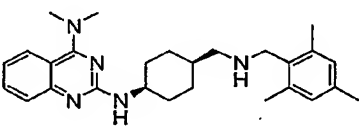
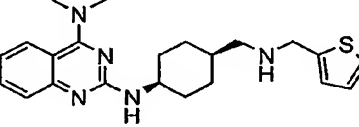
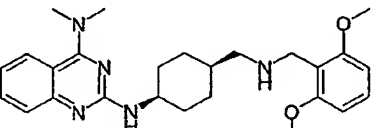
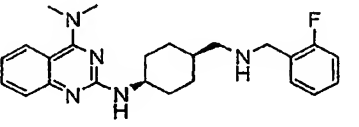
Example No.	Structure	ESI-MS	Retention Time (min)
2813	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4ccc(Br)cc4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	519.6 (M + H)	4.25
2814	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4cc(Cl)cc(Cl)c4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	509.4 (M + H)	4.49
2815	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4cc(Br)cc(OC(F)(F)F)c4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	603.0 (M + H)	4.60
2816	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4cc(C(F)(F)F)cc(C(F)(F)F)c4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	577.6 (M + H)	4.72
2817	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4c5ccccc5cc4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	491 (M + H)	4.31
2818	 <chem>Nc1nc2c(ncn2C1CCN3CC[C@H](CC3)CNS(=O)(=O)c4c5ccccc5cc4)cc3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	491.6 (M + H)	4.33

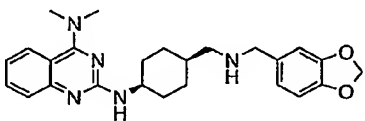
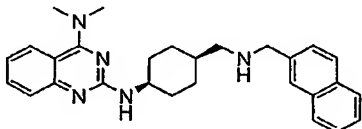
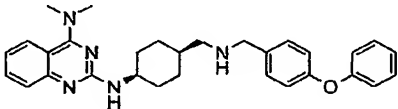
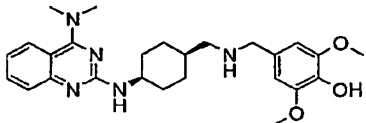
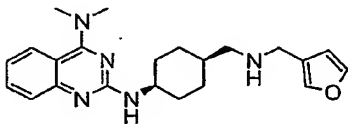
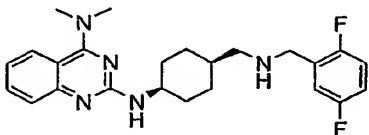
Example No.	Structure	ESI-MS	Retention Time (min)
2819	 <chem>CC1=CC=C(C=C1C2=CC=CC=C2NS(=O)(=O)NCC3CCCCC3NC4=NC5=CC=CC=C5N=C4N=CN)CC6CCCCC6</chem> $3\text{CF}_3\text{CO}_2\text{H}$	534.6 (M + H)	4.01
2820	 <chem>N=C(N)NCC1CCCCC1NS(=O)(=O)c2ccccc2</chem> 2HCl	325.4 (M + H)	3.91
2821	 <chem>N=C(N)NCC1CCCCC1NS(=O)(=O)c2ccc(Cl)cc2</chem> 2HCl	359.4 (M + H)	4.24
2822	 <chem>N=C(N)NCC1CCCCC1NS(=O)(=O)c2ccc(OC(F)(F)F)cc2</chem> 2HCl	409.4 (M + H)	4.51
2823	 <chem>N=C(N)NCC1CCCCC1NS(=O)(=O)c2ccc(C)cc2</chem> 2HCl	339.6 (M + H)	4.09
2824	 <chem>N=C(N)NCC1CCCCC1NS(=O)(=O)c2ccc(Br)cc2</chem> 2HCl	403.4 (M + H)	4.28

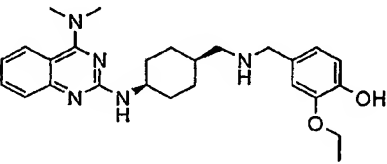
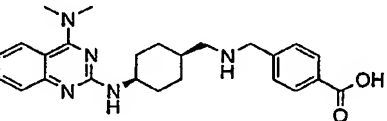
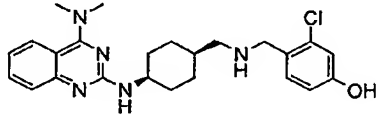
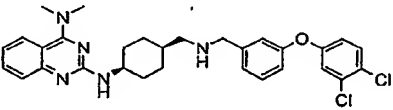
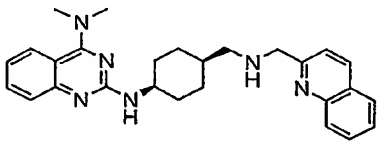
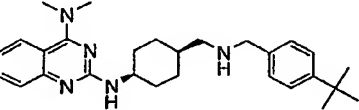
Example No.	Structure	ESI-MS	Retention Time (min)
2825	 2HCl	393.0 (M + H)	4.57
2826	 2HCl	521.6 (M + H)	4.69
2827	 2HCl	461.6 (M + H)	4.77
2828	 2HCl	375.4 (M + H)	4.33
2829	 2HCl	375.4 (M + H)	4.39
2830	 2HCl	418.8 (M + H)	4.33

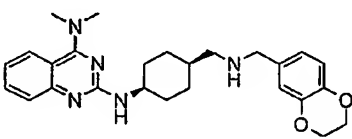
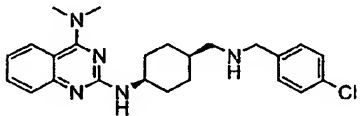
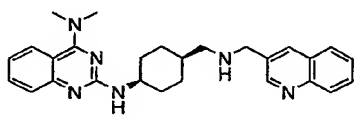
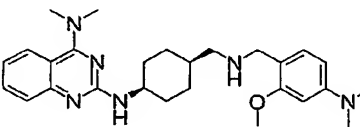
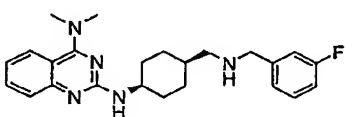
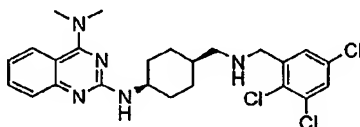
Example No.	Structure	ESI-MS	Retention Time (min)
2831	 2HCl	343.4 (M + H)	3.96
2832	 2HCl	343.4 (M + H)	4.03
2833	 2HCl	359.4 (M + H)	4.05
2834	 2HCl	359.4 (M + H)	4.24
2835	 2HCl	403.4 (M + H)	4.07
2836	 2HCl	385.4 (M + H)	4.00

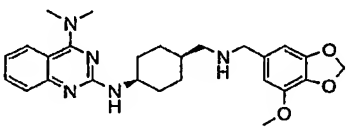
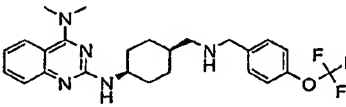
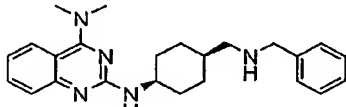
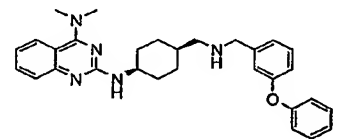
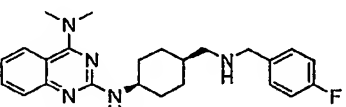
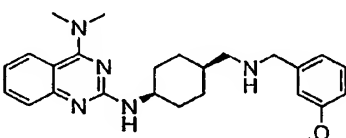
Example No.	Structure	ESI-MS	Retention Time (min)
2837	 2HCl	409.4 (M + H)	4.32
2838	 2HCl	393.6 (M + H)	4.23
2839	 2HCl	381.6 (M + H)	4.62
2840	 2HCl	330.8 (M + H)	3.83
2841	 2HCl	361.4 (M + H)	4.05
2842	 2HCl	427.4 (M + H)	4.51

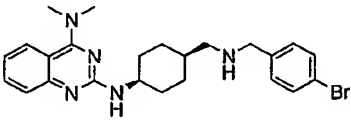
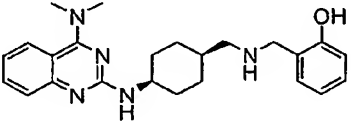
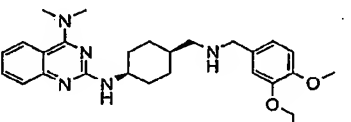
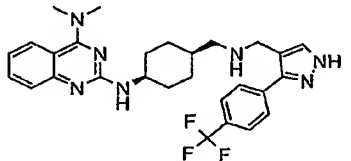
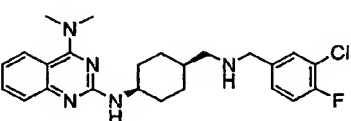
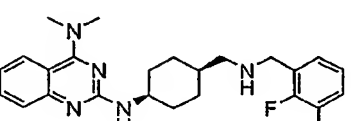
Example No.	Structure	ESI-MS	Retention Time (min)
2843	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(Cl)cc(Cl)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	458.4 (M + H)	3.22
2844	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4c[nH]c5ccccc45C#N</chem> $2\text{CF}_3\text{CO}_2\text{H}$	415.4 (M + H)	3.01
2845	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4c(C)c(C)c(C)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	432.6 (M + H)	3.26
2846	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4ccsc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	396.2 (M + H)	2.81
2847	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(OC)c(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	3.09
2848	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(F)ccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	408.4 (M + H)	2.85

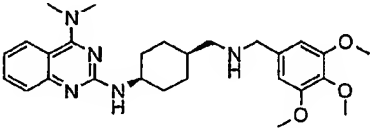
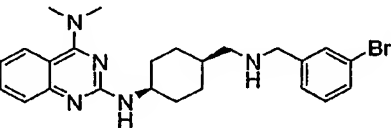
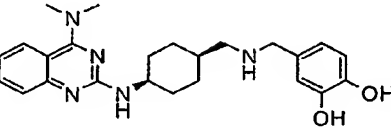
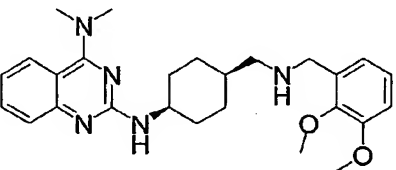
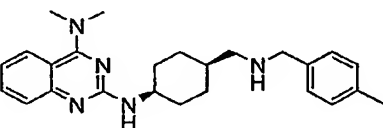
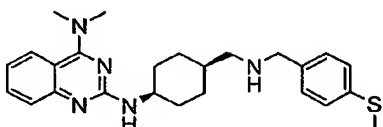
Example No.	Structure	ESI-MS	Retention Time (min)
2849	 $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	2.89
2850	 $2\text{CF}_3\text{CO}_2\text{H}$	440.0 (M + H)	3.20
2851	 $2\text{CF}_3\text{CO}_2\text{H}$	482.4 (M + H)	3.43
2852	 $2\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	2.71
2853	 $2\text{CF}_3\text{CO}_2\text{H}$	380.2 (M + H)	2.72
2854	 $2\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	2.91

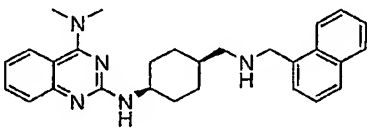
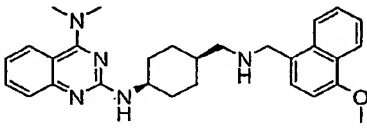
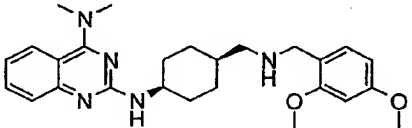
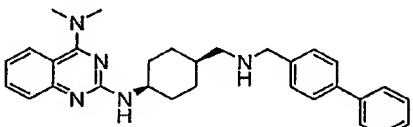
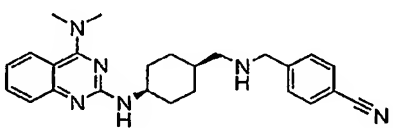
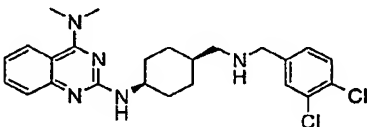
Example No.	Structure	ESI-MS	Retention Time (min)
2855	 $2\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	2.82
2856	 $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	2.69
2857	 $2\text{CF}_3\text{CO}_2\text{H}$	440.0 (M + H)	2.85
2858	 $2\text{CF}_3\text{CO}_2\text{H}$	550.6 (M + H)	3.80
2859	 $3\text{CF}_3\text{CO}_2\text{H}$	441.4 (M + H)	3.03
2860	 $2\text{CF}_3\text{CO}_2\text{H}$	446.6 (M + H)	3.41

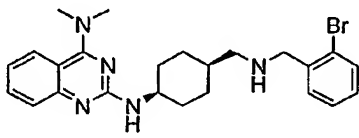
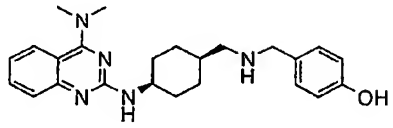
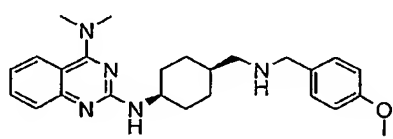
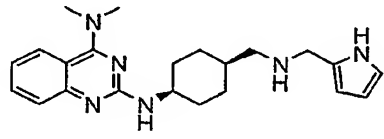
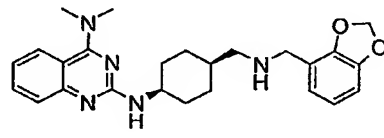
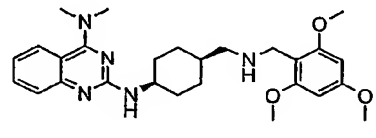
Example No.	Structure	ESI-MS	Retention Time (min)
2861	 $2\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	2.91
2862	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	3.05
2863	 $3\text{CF}_3\text{CO}_2\text{H}$	441.4 (M + H)	2.68
2864	 $3\text{CF}_3\text{CO}_2\text{H}$	463.4 (M + H)	2.76
2865	 $2\text{CF}_3\text{CO}_2\text{H}$	408.4 (M + H)	2.91
2866	 $2\text{CF}_3\text{CO}_2\text{H}$	492.2 (M + H)	3.30

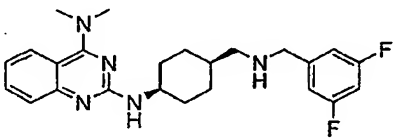
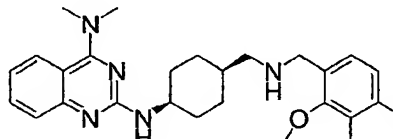
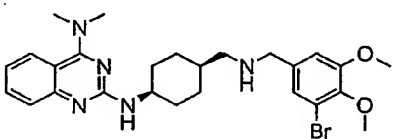
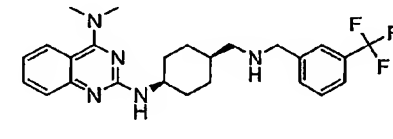
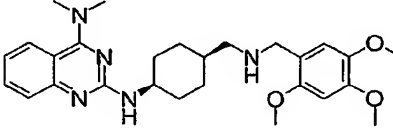
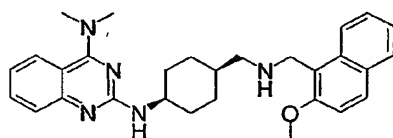
Example No.	Structure	ESI-MS	Retention Time (min)
2867	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>COc1ccc2c(c1)OCO2</chem> <chem>2CF3CO2H</chem>	464.2 (M + H)	2.93
2868	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>COc1ccc(OC(F)(F)F)cc1</chem> <chem>2CF3CO2H</chem>	474.4 (M + H)	3.27
2869	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>c1ccccc1</chem> <chem>2CF3CO2H</chem>	390.6 (M + H)	2.88
2870	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>c1ccc(cc1)Oc2ccccc2</chem> <chem>2CF3CO2H</chem>	482.2 (M + H)	3.43
2871	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>Fc1ccc(cc1)</chem> <chem>2CF3CO2H</chem>	408.4 (M + H)	2.91
2872	 <chem>CN(C)c1nc2c(ncn2C1)C3=CC=CC=C3</chem> <chem>COc1ccc(cc1)</chem> <chem>2CF3CO2H</chem>	420.4 (M + H)	2.91

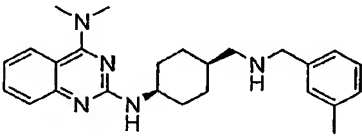
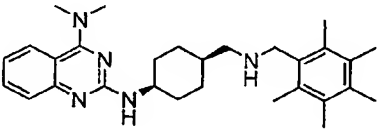
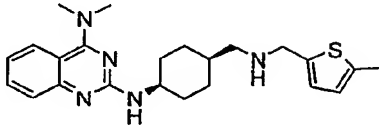
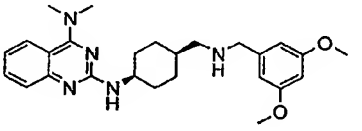
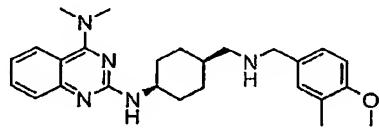
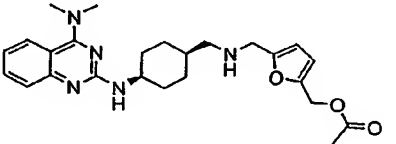
Example No.	Structure	ESI-MS	Retention Time (min)
2873	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC=CC=C4N3)C5=CC=CC=C5Br</chem> $2\text{CF}_3\text{CO}_2\text{H}$	468.2 (M + H)	3.09
2874	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC(=C(C=C4)O)N3)C5=CC=CC=C5O</chem> $2\text{CF}_3\text{CO}_2\text{H}$	406.4 (M + H)	2.80
2875	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC(=C(C=C4)OCC)N3)C5=CC=CC=C5OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	464.2 (M + H)	2.97
2876	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC(=C(C=C4)C(F)(F)F)N3)C5=CC=CC=C5C(F)(F)F</chem> $3\text{CF}_3\text{CO}_2\text{H}$	524.6 (M + H)	3.12
2877	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC(=C(C=C4)Cl)N3)C5=CC(=C(C=C5)F)N6=CC=CC=C6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	442.4 (M + H)	3.10
2878	 <chem>CN1C=NC2=C(N1)C=CC=C2N3C(=NC4=CC(=C(C=C4)F)N3)C5=CC(=C(C=C5)F)N6=CC=CC=C6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	2.90

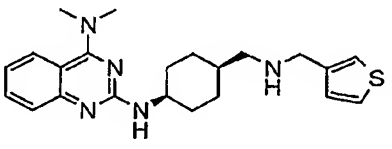
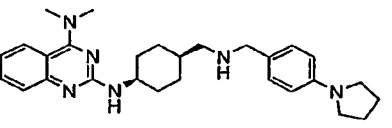
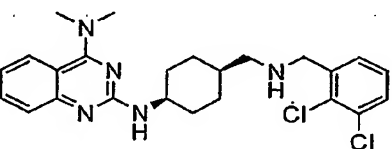
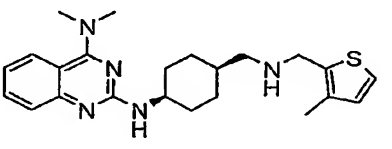
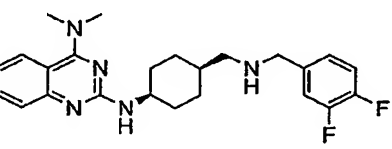
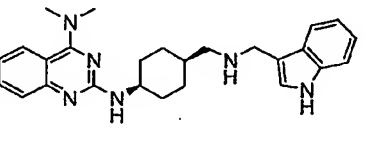
Example No.	Structure	ESI-MS	Retention Time (min)
2879	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4cc(OC)c(OC)c(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	2.89
2880	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(Br)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	468.2 (M + H)	3.07
2881	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(O)c(O)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	422.4 (M + H)	2.61
2882	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4cc(OC)c(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	2.93
2883	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	404.6 (M + H)	3.01
2884	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(I)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	3.08

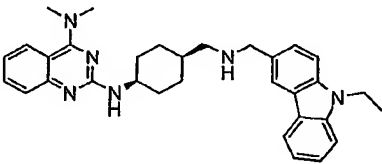
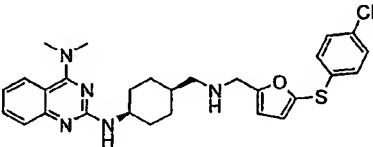
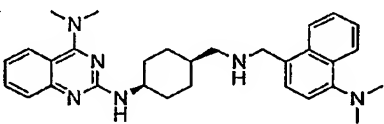
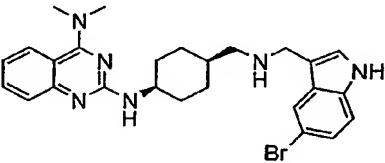
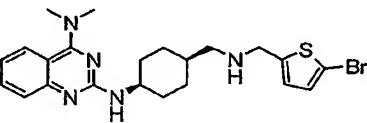
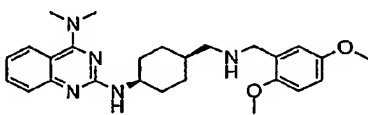
Example No.	Structure	ESI-MS	Retention Time (min)
2885	 2CF ₃ CO ₂ H	440.0 (M + H)	3.18
2886	 2CF ₃ CO ₂ H	470.4 (M + H)	3.25
2887	 2CF ₃ CO ₂ H	450.0 (M + H)	3.01
2888	 2CF ₃ CO ₂ H	466.4 (M + H)	3.40
2889	 2CF ₃ CO ₂ H	415.4 (M + H)	2.83
2890	 2CF ₃ CO ₂ H	458.4 (M + H)	3.25

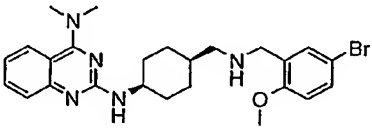
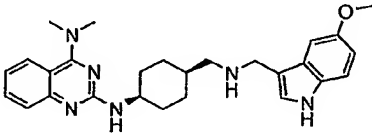
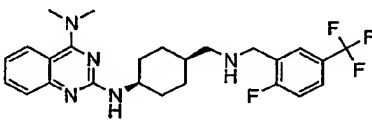
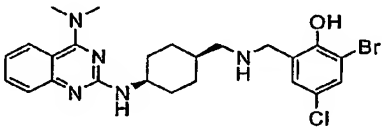
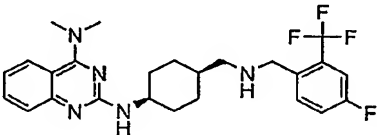
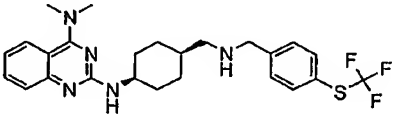
Example No.	Structure	ESI-MS	Retention Time (min)
2891	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4cc(Br)ccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	468.2 (M + H)	3.00
2892	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(O)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	406.4 (M + H)	2.66
2893	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.92
2894	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4c[nH]c5ccccc45</chem> $3\text{CF}_3\text{CO}_2\text{H}$	379.4 (M + H)	2.71
2895	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4ccc5c(c4)OCO5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	2.87
2896	 <chem>CN(C)c1nc2c(ncnc2c1)N[C@H]3CCCC[C@H]3CNCCc4c(OC)c(OC)c(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	3.17

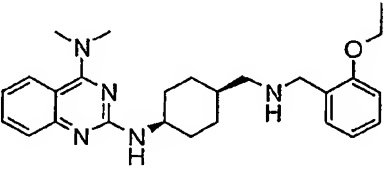
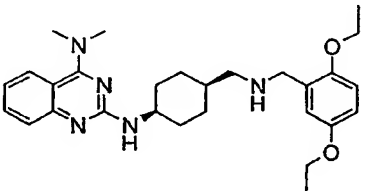
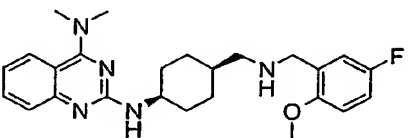
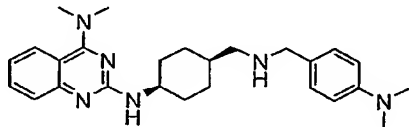
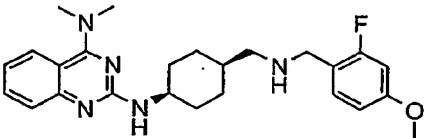
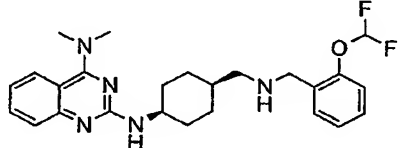
Example No.	Structure	ESI-MS	Retention Time (min)
2897	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1ccc(F)c(F)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	2.98
2898	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1cc(OC)c(OC)c(OC)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	2.99
2899	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1cc(OC)c(Br)c(OC)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	528.4 (M + H)	3.15
2900	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1ccc(C(F)(F)F)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	458.4 (M + H)	3.19
2901	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1cc(OC)c(OC)c(OC)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	2.92
2902	 <chem>CN(C)c1nc2c(ncnc12)N[C@H]3CCCC[C@H]3CNCCc1ccc2cc(OC)ccc2c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	470.4 (M + H)	3.27

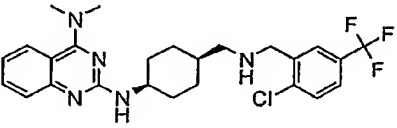
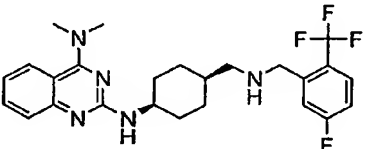
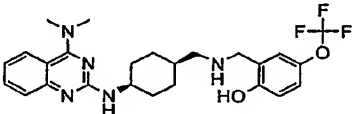
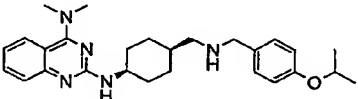
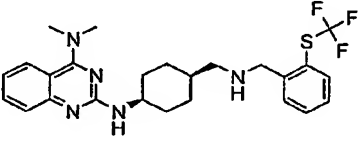
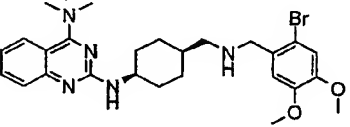
Example No.	Structure	ESI-MS	Retention Time (min)
2903	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4ccc(C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	404.6 (M + H)	2.87
2904	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4c(C)c(C)c(C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	3.48
2905	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(C)s4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	410.4 (M + H)	2.96
2906	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(OC)c(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	3.03
2907	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc(OC)ccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.08
2908	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CNc4cc5oc(cc5O4)COC(=O)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	452.2 (M + H)	2.79

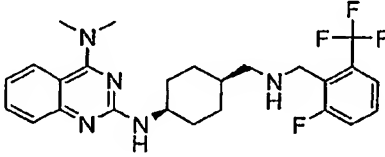
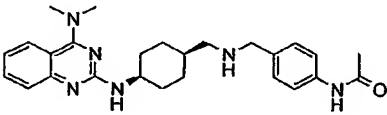
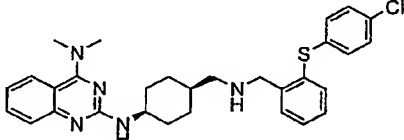
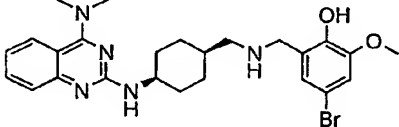
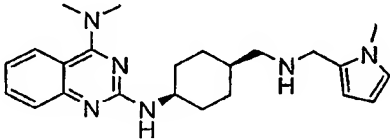
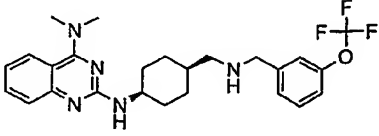
Example No.	Structure	ESI-MS	Retention Time (min)
2909	 $2\text{CF}_3\text{CO}_2\text{H}$	396.2 (M + H) /	2.81
2910	 $3\text{CF}_3\text{CO}_2\text{H}$	459.4 (M + H)	3.21
2911	 $2\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.08
2912	 $2\text{CF}_3\text{CO}_2\text{H}$	410.4 (M + H)	2.88
2913	 $2\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	3.01
2914	 $3\text{CF}_3\text{CO}_2\text{H}$	429.4 (M + H)	2.97

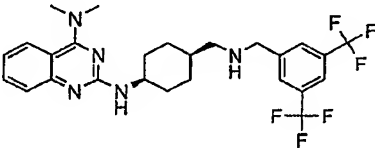
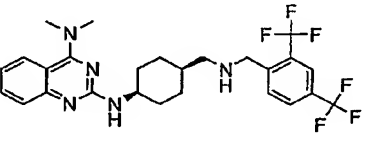
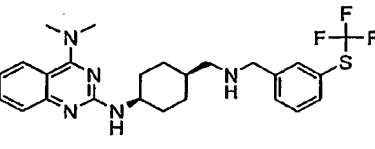
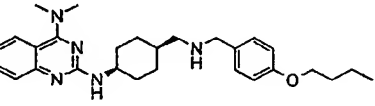
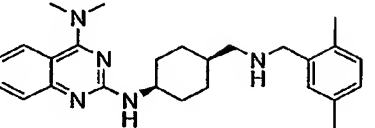
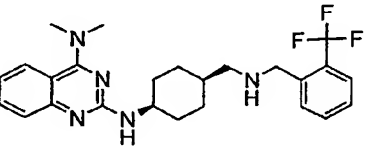
Example No.	Structure	ESI-MS	Retention Time (min)
2915	 <chem>CC1=CC=C2C(=C1)C(=NN2)C(=N)N[C@H]3CCCC[C@H]3CNCC4=CC=CC=C5C4=CC=CC5N(C)C</chem> $3\text{CF}_3\text{CO}_2\text{H}$	507.2 (M + H)	3.53
2916	 <chem>CC1=CC=C2C(=C1)C(=N)N=C2N[C@H]3CCCC[C@H]3CNCC4=CC=C(C=C4)SC5=CC=CC=C5Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	522.4 (M + H)	3.56
2917	 <chem>CC1=CC=C2C(=C1)C(=N)N=C2N[C@H]3CCCC[C@H]3CNCC4=CC=CC=C5C4=CC=CC5N(C)C</chem> $3\text{CF}_3\text{CO}_2\text{H}$	483.2 (M + H)	2.80
2918	 <chem>CC1=CC=C2C(=C1)C(=N)N=C2N[C@H]3CCCC[C@H]3CNCC4=CC=CC=C5C4=CC=CC5N(C)C</chem> $3\text{CF}_3\text{CO}_2\text{H}$	507.2 (M + H)	3.27
2919	 <chem>CC1=CC=C2C(=C1)C(=N)N=C2N[C@H]3CCCC[C@H]3CNCC4=CC=C(C=C4)SC5=CC=CC=C5Br</chem> $2\text{CF}_3\text{CO}_2\text{H}$	474.2 (M + H)	3.10
2920	 <chem>CC1=CC=C2C(=C1)C(=N)N=C2N[C@H]3CCCC[C@H]3CNCC4=CC=C(C=C4)SC5=CC=CC=C5OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	3.00

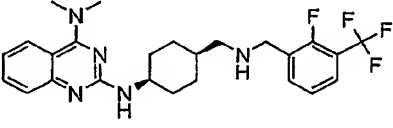
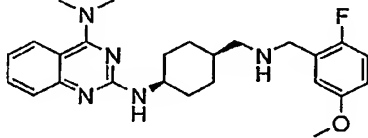
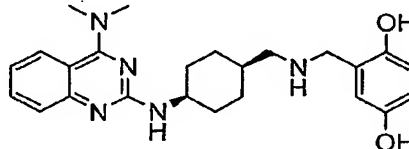
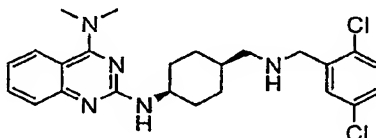
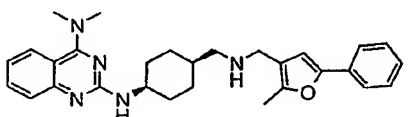
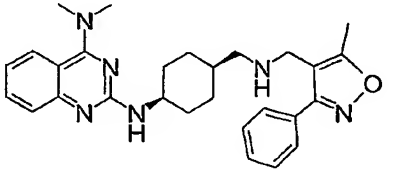
Example No.	Structure	ESI-MS	Retention Time (min)
2921	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CC(OC)=CC4Br</chem> $2\text{CF}_3\text{CO}_2\text{H}$	498.4 (M + H)	3.15
2922	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CC(OC)=CC4Cc5c[nH]c6ccccc56</chem> $3\text{CF}_3\text{CO}_2\text{H}$	459.4 (M + H)	2.99
2923	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CC(F)=CC4C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.0 (M + H)	3.10
2924	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CC(O)=CC4BrCl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	518.2 (M + H)	3.10
2925	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CC(F)=CC4C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.12
2926	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3NC4C=CCSC4C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.35

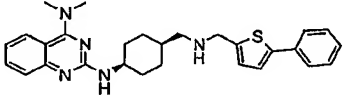
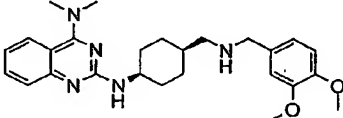
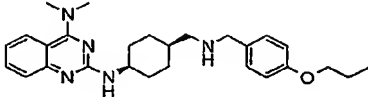
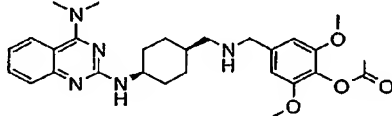
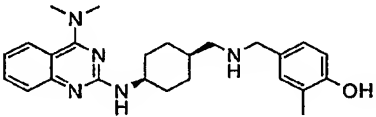
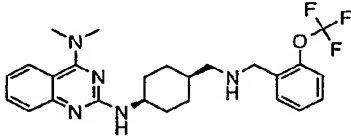
Exemplé No.	Structure	ESI-MS	Retention Time (min)
2927	 2CF ₃ CO ₂ H	434.4 (M + H)	3.11
2928	 2CF ₃ CO ₂ H	478.4 (M + H)	3.29
2929	 2CF ₃ CO ₂ H	438.2 (M + H)	3.01
2930	 3CF ₃ CO ₂ H	433.4 (M + H)	2.59
2931	 2CF ₃ CO ₂ H	438.2 (M + H)	2.90
2932	 2CF ₃ CO ₂ H	456.2 (M + H)	3.10

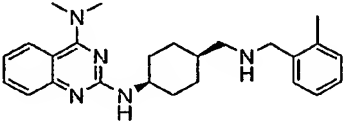
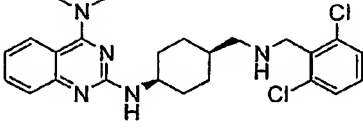
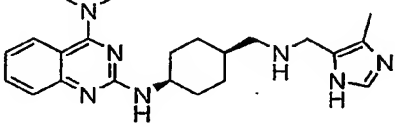
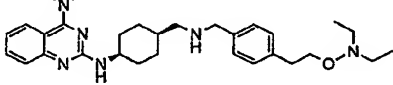
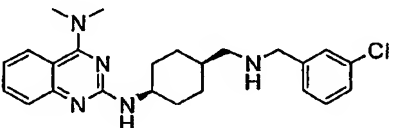
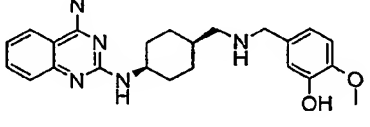
Example No.	Structure	ESI-MS	Retention Time (min)
2933	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3CN[C@@H]4C=CC(=CC=C4)C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.2 (M + H)	3.25
2934	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3CN[C@@H]4C=CC(=CC=C4)C(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.11
2935	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3CN[C@@H]4C=CC(=CC=C4)C(F)(F)OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.20
2936	 <chem>CC(C)Oc1ccc(cc1)CN[C@@H]2CCCC[C@H]2N[C@H]3C=NC4=CC=CC=C4N(C)C3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.17
2937	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3CN[C@@H]4C=CC(=CC=C4)SC(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	489.6 (M + H)	3.31
2938	 <chem>COc1cc(OC)c(Br)cc1CN[C@@H]2CCCC[C@H]2N[C@H]3C=NC4=CC=CC=C4N(C)C3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	528.2 (M + H)	3.03

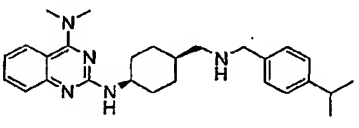
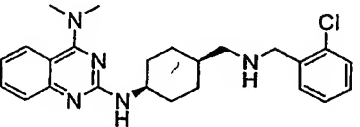
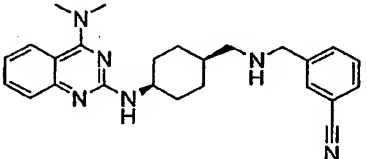
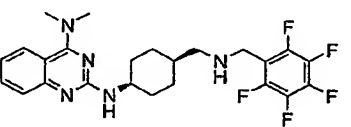
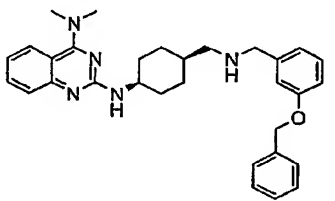
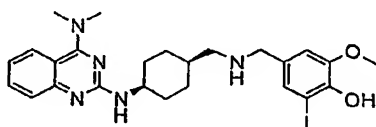
Example No.	Structure	ESI-MS	Retention Time (min)
2939	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)C(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	2.99
2940	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)NC(=O)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	447.4 (M + H)	2.66
2941	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)SC5=CC=C(Cl)C=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	532.4 (M + H)	3.66
2942	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)C(O)C(OC)C(Br)=C4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	514.4 (M + H)	3.08
2943	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)C5=CC=CN5</chem> $3\text{CF}_3\text{CO}_2\text{H}$	393.4 (M + H)	2.79
2944	 <chem>CN(C)c1nc2ccccc2n1NC3CCCCC3CCNC4C=CC(=CC=C4)C5=CC=C(OC(F)(F)F)C=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	3.24

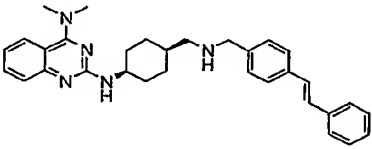
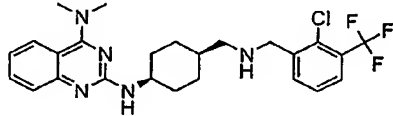
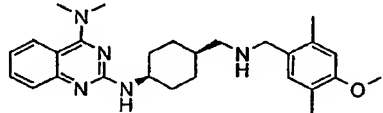
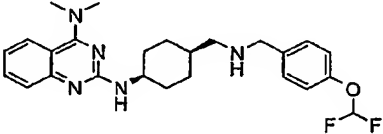
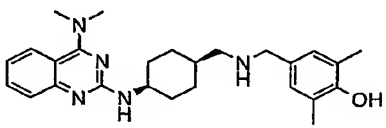
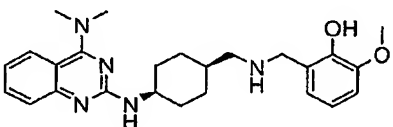
Example No.	Structure	ESI-MS	Retention Time (min)
2945	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(C(F)(F)F)C(F)(F)F=C5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	526.6 (M + H)	3.44
2946	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(C(F)(F)F)C(F)(F)F=C5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	526.6 (M + H)	3.42
2947	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(C(F)(F)F)S(=O)(=O)N5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.35
2948	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(COCC)C=C5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.43
2949	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(C)C=C5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	3.13
2950	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4CCCCC4NC5=CC=C(C(F)(F)F)C=C5)c3ccccc13</chem> $2\text{CF}_3\text{CO}_2\text{H}$	458.4 (M + H)	3.10

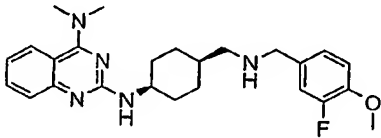
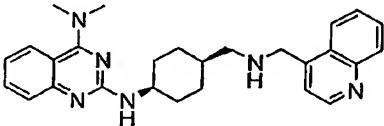
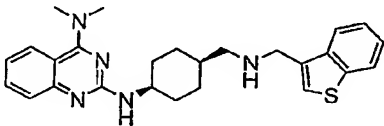
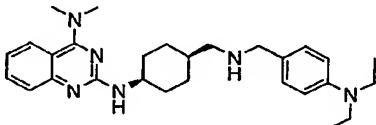
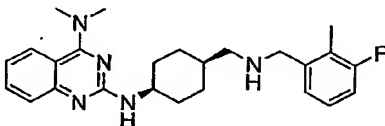
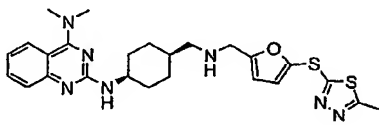
Example No.	Structure	ESI-MS	Retention Time (min)
2951	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC=C(C(F)(F)F)C=C4)C3=CC=CC=C3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	3.19
2952	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC=C(F)C=C4)C3=CC=CC=C3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	2.95
2953	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC(=C(C=C4)O)O</chem> $2\text{CF}_3\text{CO}_2\text{H}$	422.4 (M + H)	2.61
2954	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC(=C(C=C4)Cl)Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.07
2955	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC5=C(C=C4)OC(=C5)c6ccccc6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	470.4 (M + H)	3.45
2956	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCC4=CC5=C(C=C4)ONC5C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	471.6 (M + H)	2.88

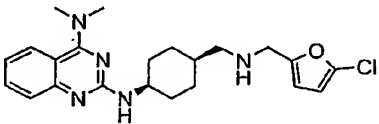
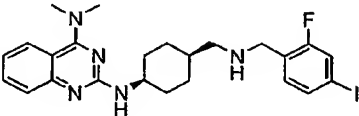
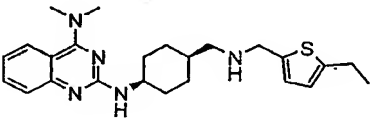
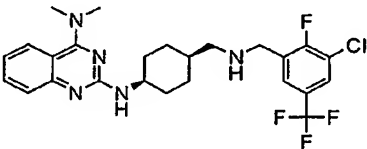
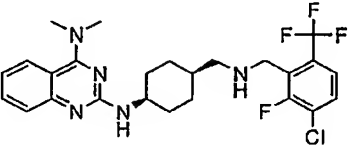
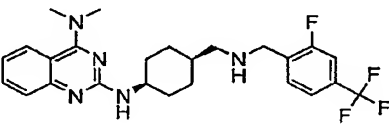
Example No.	Structure	ESI-MS	Retention Time (min)
2957	 $2\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	3.36
2958	 $2\text{CF}_3\text{CO}_2\text{H}$	450 (M + H)	2.75
2959	 $2\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.20
2960	 $2\text{CF}_3\text{CO}_2\text{H}$	508.4 (M + H)	3.00
2961	 $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.80
2962	 $2\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	3.20

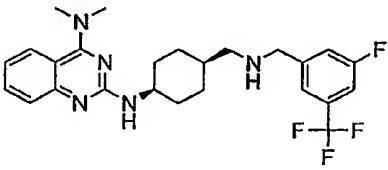
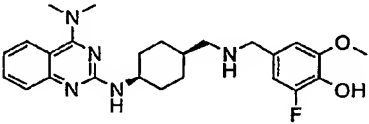
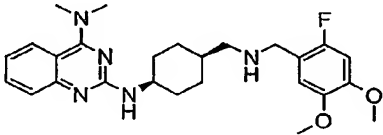
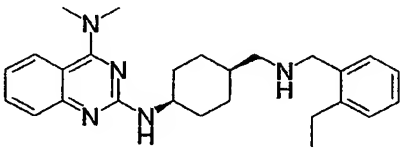
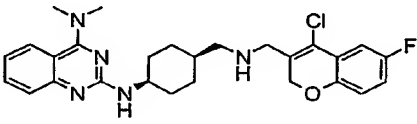
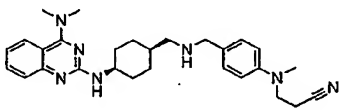
Example No.	Structure	ESI-MS	Retention Time (min)
2963	 $2\text{CF}_3\text{CO}_2\text{H}$	404.4 (M + H)	2.87
2964	 $2\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.00
2965	 $3\text{CF}_3\text{CO}_2\text{H}$	394.4 (M + H)	2.30
2966	 $2\text{CF}_3\text{CO}_2\text{H}$	505.4 (M + H)	2.60
2967	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	3.00
2968	 $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	2.71

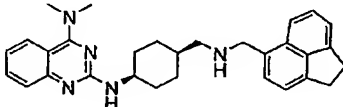
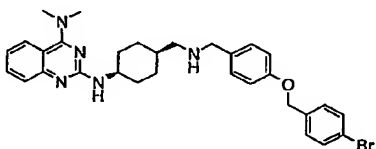
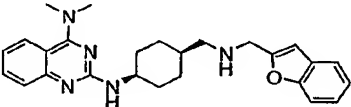
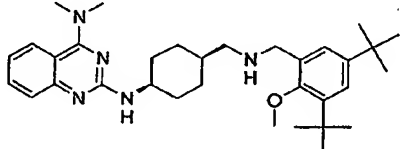
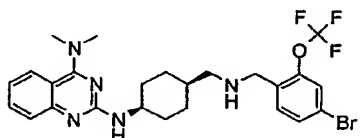
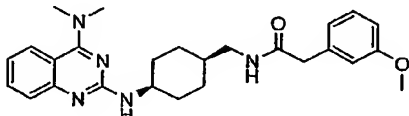
Example No.	Structure	ESI-MS	Retention Time (min)
2969	 <chem>CC(C)c1ccc(cc1)CNCC2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	432.4 (M + H)	3.30
2970	 <chem>Clc1ccc(cc1)CNCC2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	2.95
2971	 <chem>N#Cc1ccc(cc1)CNCC2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	415.4 (M + H)	2.79
2972	 <chem>Fc1cc(F)c(F)c(F)c1CNCC2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	3.00
2973	 <chem>c1ccc(cc1)OCCc2ccc(cc2)CNCC3CCCCC3Nc4nc5ccccc5n4C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	496.2 (M + H)	3.46
2974	 <chem>COc1cc(O)ccc1CNCC2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	562.2 (M + H)	2.99

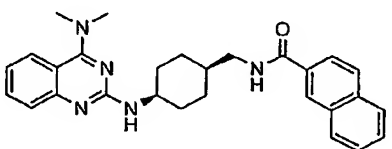
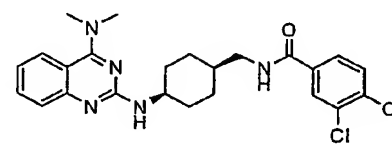
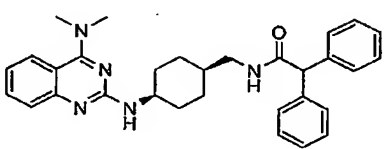
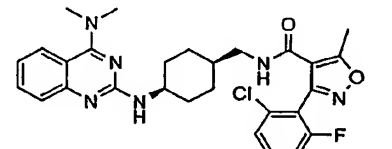
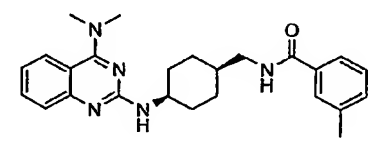
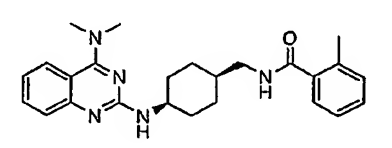
Example No.	Structure	ESI-MS	Retention Time (min)
2975	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(cc4)/C=C/c5ccccc5)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.4 (M + H)	3.64
2976	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(cc4)C(F)(F)F)Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.2 (M + H)	3.25
2977	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(OC)cc4)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.22
2978	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(OC(F)(F)F)cc4)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	456.2 (M + H)	3.09
2979	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(O)cc4)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	2.89
2980	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3CNCCc4ccc(OC)cc4)O</chem> $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	2.79

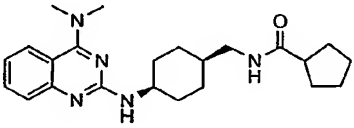
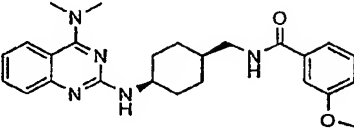
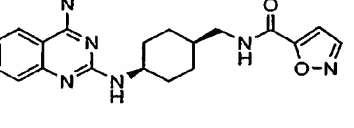
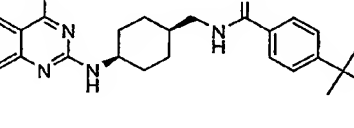
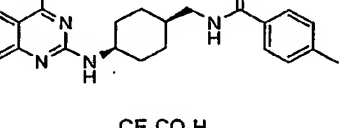
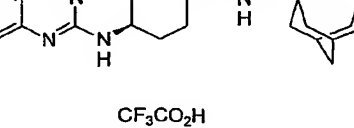
Example No.	Structure	ESI-MS	Retention Time (min)
2981	 $2\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	2.91
2982	 $3\text{CF}_3\text{CO}_2\text{H}$	441.4 (M + H)	2.55
2983	 $2\text{CF}_3\text{CO}_2\text{H}$	446.4 (M + H)	3.13
2984	 $3\text{CF}_3\text{CO}_2\text{H}$	461.4 (M + H)	2.46
2985	 $2\text{CF}_3\text{CO}_2\text{H}$	422.2 (M + H)	3.01
2986	 $2\text{CF}_3\text{CO}_2\text{H}$	510.2 (M + H)	2.85

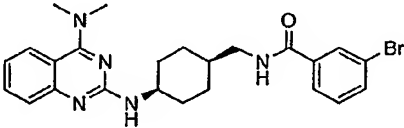
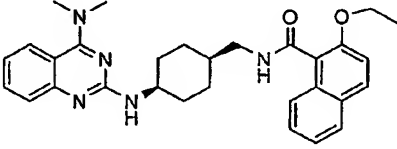
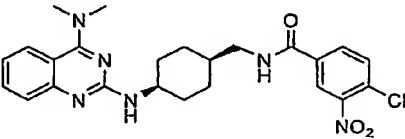
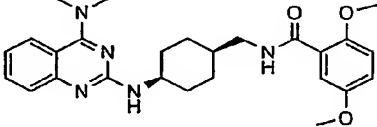
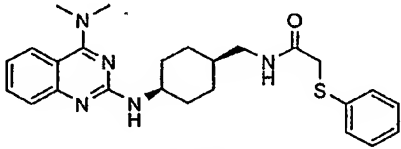
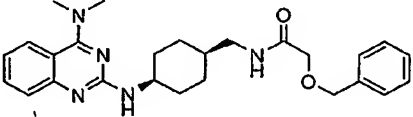
Example No.	Structure	ESI-MS	Retention Time (min)
2987	 $2\text{CF}_3\text{CO}_2\text{H}$	414.4 (M + H)	2.86
2988	 $2\text{CF}_3\text{CO}_2\text{H}$	534.2 (M + H)	3.13
2989	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	3.08
2990	 $2\text{CF}_3\text{CO}_2\text{H}$	510.4 (M + H)	3.32
2991	 $2\text{CF}_3\text{CO}_2\text{H}$	510.4 (M + H)	3.17
2992	 $2\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	3.17

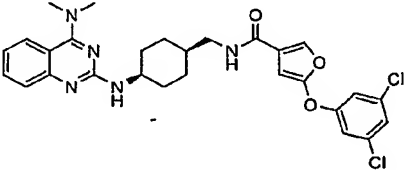
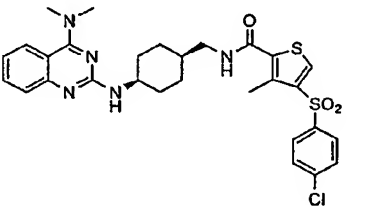
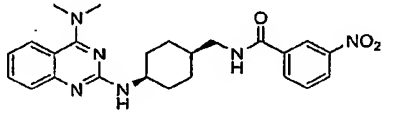
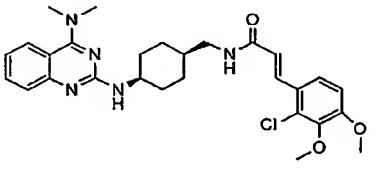
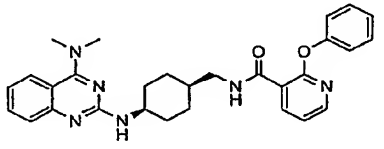
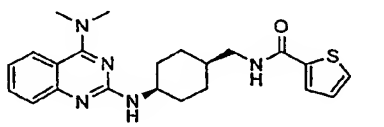
Example No.	Structure	ESI-MS	Retention Time (min)
2993	 $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.21
2994	 $2\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	2.77
2995	 $2\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	2.89
2996	 $2\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	3.12
2997	 $2\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	3.29
2998	 $3\text{CF}_3\text{CO}_2\text{H}$	472.6 (M + H)	2.99

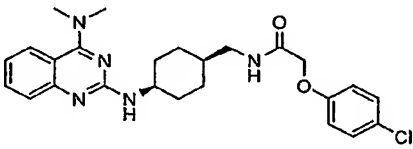
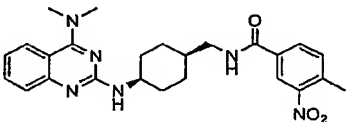
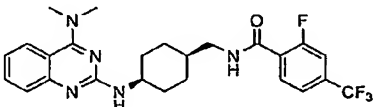
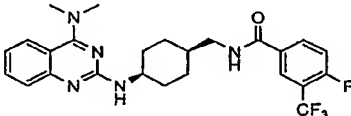
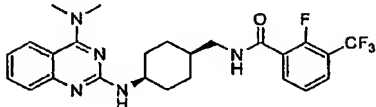
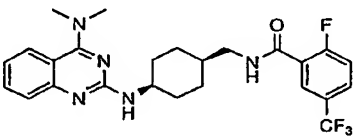
Example No.	Structure	ESI-MS	Retention Time (min)
2999	 $2\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	3.37
3000	 $2\text{CF}_3\text{CO}_2\text{H}$	574.2 (M + H)	3.64
3001	 $2\text{CF}_3\text{CO}_2\text{H}$	430.4 (M + H)	3.05
3002	 $2\text{CF}_3\text{CO}_2\text{H}$	532.4 (M + H)	4.05
3003	 $2\text{CF}_3\text{CO}_2\text{H}$	552.0 (M + H)	3.37
3004	 $\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.51

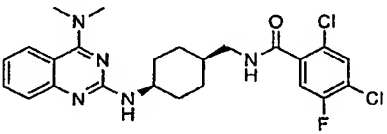
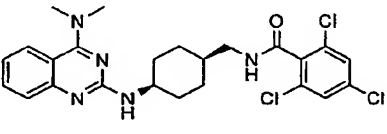
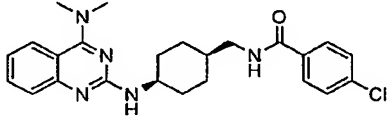
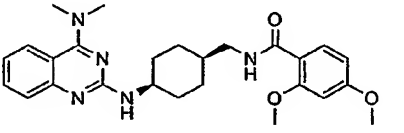
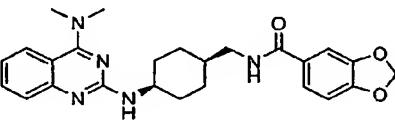
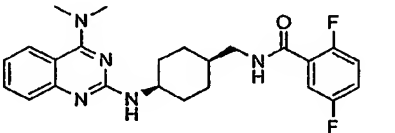
Example No.	Structure	ESI-MS	Retention Time (min)
3005	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)Nc5ccc6ccccc6c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	3.91
3006	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)Nc5ccc(Cl)c(Cl)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	4.02
3007	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)N(C(=O)c5ccccc5)c6ccccc6</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.4 (M + H)	4.01
3008	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)Nc5c(Cl)c6ccccc6n5C(F)=O</chem> $\text{CF}_3\text{CO}_2\text{H}$	537.4 (M + H)	3.77
3009	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)Nc5ccc(C)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	3.63
3010	 <chem>CC1=CN2C(=N1)c3ccccc3N2C4CCCCC4CC(=O)Nc5cccc(C)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	3.51

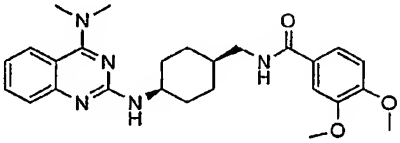
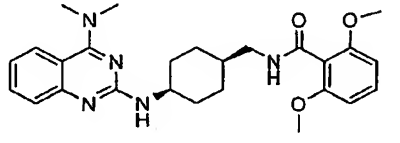
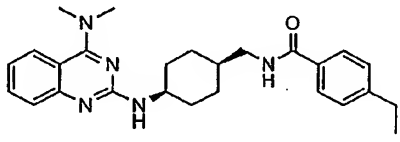
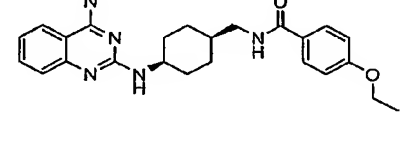
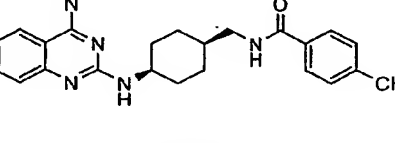
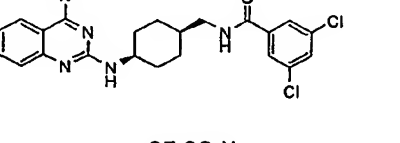
Example No.	Structure	ESI-MS	Retention Time (min)
3011	 CF ₃ CO ₂ H	396.2 (M + H)	3.47
3012	 CF ₃ CO ₂ H	434.4 (M + H)	3.52
3013	 CF ₃ CO ₂ H	395.4 (M + H)	3.15
3014	 CF ₃ CO ₂ H	460.2 (M + H)	4.03
3015	 CF ₃ CO ₂ H	418.6 (M + H)	3.65
3016	 CF ₃ CO ₂ H	462.2 (M + H)	4.09

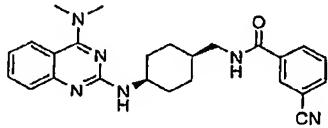
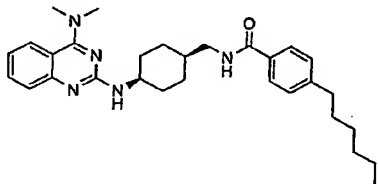
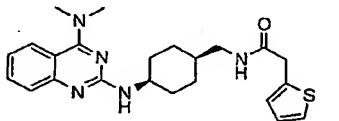
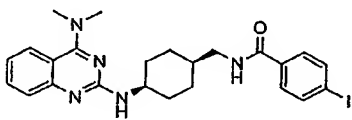
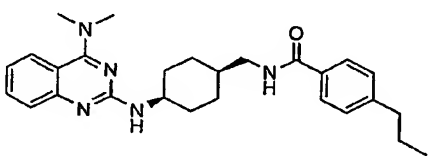
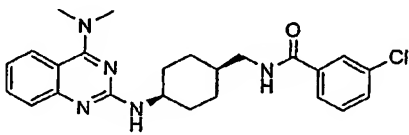
Example No.	Structure	ESI-MS	Retention Time (min)
3017	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=CC=C4C(=O)N3C5=CC=C(C=C5)Br</chem> $\text{CF}_3\text{CO}_2\text{H}$	484.2 (M + H)	3.79
3018	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=CC=C4C(=O)N3C5=CC=C(C=C5)OCC</chem> $\text{CF}_3\text{CO}_2\text{H}$	498.6 (M + H)	3.88
3019	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=C(C=C4)C(=O)N3C5=CC=C(C=C5)[N+](=O)[O-]Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	483.2 (M + H)	3.80
3020	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=C(C=C4)C(=O)N3C5=CC=C(C=C5)OC</chem> $\text{CF}_3\text{CO}_2\text{H}$	478.2 (M + H)	3.49
3021	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=CC=C4C(=O)N3C5=CC=C(C=C5)SC6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.0 (M + H)	3.61
3022	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3=CC=CC=C3N3C4=CC=CC=C4C(=O)N3C5=CC=C(C=C5)OCC6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	448.2 (M + H)	3.70

Example No.	Structure	ESI-MS	Retention Time (min)
3023	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)Nc5oc(cc5Oc6cc(Cl)cc(Cl)c6)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	554.4 (M + H)	4.41
3024	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)Nc5sc(Cc6ccc(Cl)cc6S(=O)(=O)c7ccccc7Cl)c5C</chem> $\text{CF}_3\text{CO}_2\text{H}$	598.2 (M + H)	4.03
3025	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)Nc5ccc(cc5)[N+](=O)[O-]</chem> $\text{CF}_3\text{CO}_2\text{H}$	499.2 (M + H)	3.59
3026	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)N/C=C/c5cc(OC)c(Cl)c(OC)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	524.6 (M + H)	3.84
3027	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)Nc5ccc6ccccc6o5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	497.4 (M + H)	3.80
3028	 <chem>CC1=NC2=CC=CC=C2N1N(C)C3=CC=CC=C3N3C4CCCCC4CC(=O)Nc5ccsc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	410.2 (M + H)	3.43

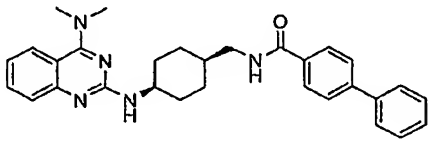
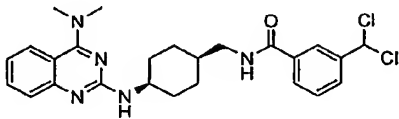
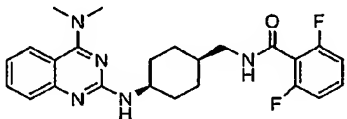
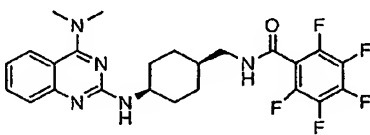
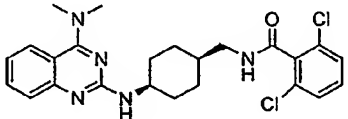
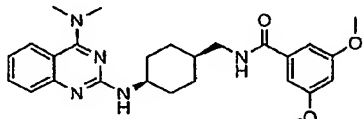
Example No.	Structure	ESI-MS	Retention Time (min)
3029	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)COc4ccc(Cl)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	468.2 (M + H)	3.77
3030	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)Cc4ccc([N+](=O)[O-])cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	463.2 (M + H)	3.73
3031	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)Cc4cc(F)c(C(F)(F)F)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.91
3032	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)Cc4cc(F)c(C(F)(F)F)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.94
3033	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)Cc4cc(F)c(C(F)(F)F)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.85
3034	 <chem>CC1=NC2=CC=CC=C2N(C)C1=NC(NC3CCCCC3CNC(=O)Cc4cc(F)c(C(F)(F)F)cc4)C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.87

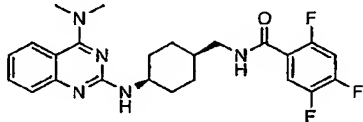
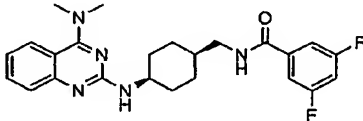
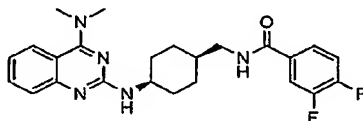
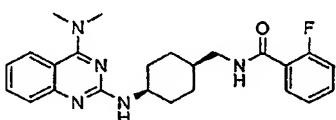
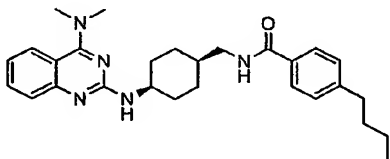
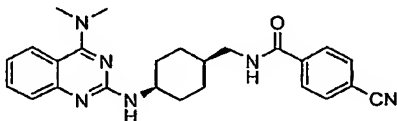
Example No.	Structure	ESI-MS	Retention Time (min)
3041	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3cc(Cl)cc(F)c3Cl)</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.4 (M + H)	3.82
3042	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3cc(Cl)cc(Cl)c3Cl)</chem> $\text{CF}_3\text{CO}_2\text{H}$	508.0 (M + H)	3.85
3043	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3ccc(Cl)cc3)</chem> $\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	3.71
3044	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3cc(OC)cc(OC)c3)</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.2 (M + H)	3.65
3045	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3cc4occc4cc3)</chem> $\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.47
3046	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN(C(=O)c3cc(F)cc(F)c3)</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	3.59

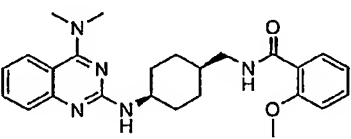
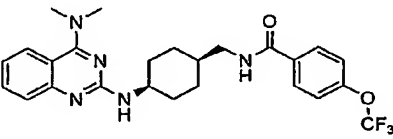
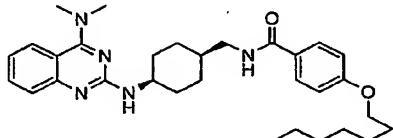
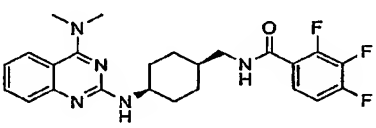
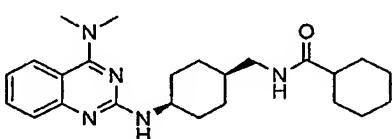
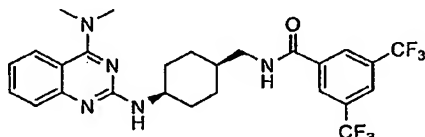
Example No.	Structure	ESI-MS	Retention Time (min)
3047	 <chem>COc1cc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)c(OC)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.2 (M + H)	3.36
3048	 <chem>COc1cc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)c(OC)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	3.39
3049	 <chem>CCc1ccc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	432.4 (M + H)	3.81
3050	 <chem>COc1ccc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	3.69
3051	 <chem>Clc1ccc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	3.69
3052	 <chem>Clc1cc(NC(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)cc(Cl)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	4.03

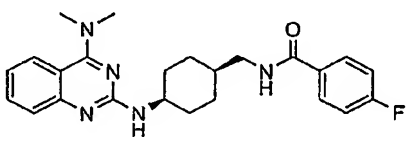
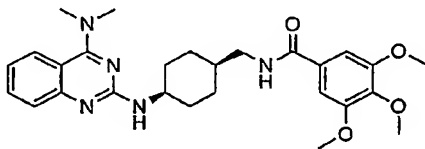
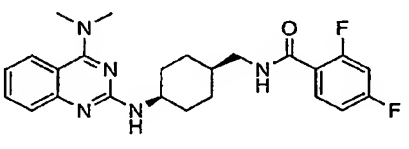
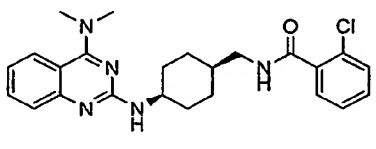
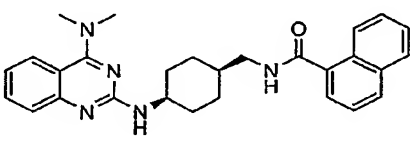
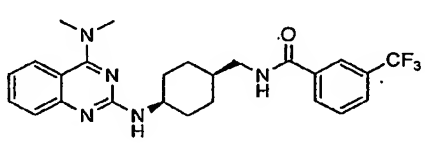
Example No.	Structure	ESI-MS	Retention Time (min)
3053	 <chem>CC1=NC2=CC=CC=C2N(C)N1NC(=O)c3ccc(C#N)cc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	429.2 (M + H)	3.47
3054	 <chem>CCCCCc1ccc(cc1)C(=O)NC2CCC(CC2)Nc3nc4c(ncn3C)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	4.60
3055	 <chem>Cc1cc2c(ncn1C)ccccc2NC(=O)Cc3ccsc3</chem> $\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	3.41
3056	 <chem>CIc1ccc(cc1)C(=O)NC2CCC(CC2)Nc3nc4c(ncn3C)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	530.2 (M + H)	3.83
3057	 <chem>CCCc1ccc(cc1)C(=O)NC2CCC(CC2)Nc3nc4c(ncn3C)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	446.4 (M + H)	4.02
3058	 <chem>Clc1ccc(cc1)C(=O)NC2CCC(CC2)Nc3nc4c(ncn3C)ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	3.70

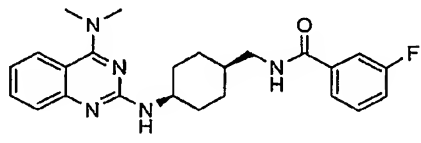
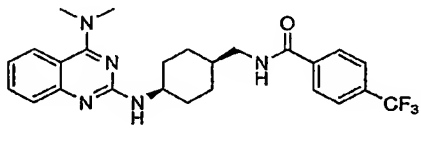
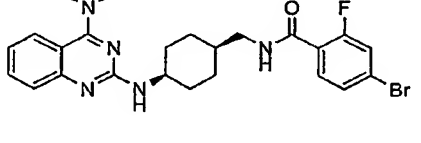
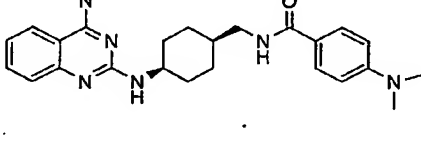
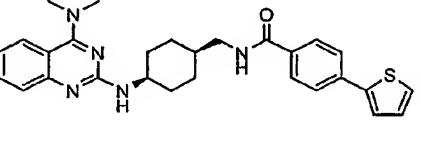
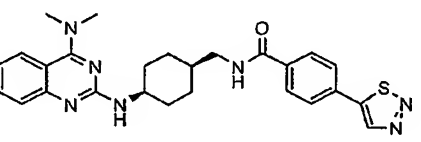
[illegible]

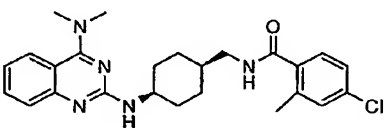
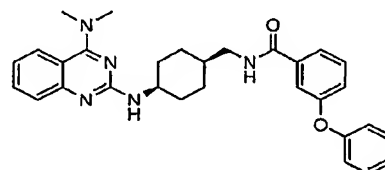
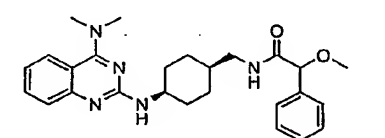
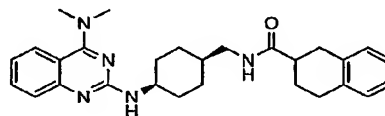
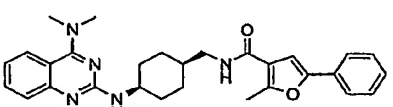
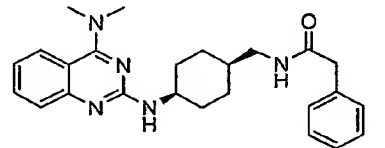
Example No.	Structure	ESI-MS	Retention Time (min)
3065	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=CC=CC=C5</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	4.09
3066	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=C(C=C4)C(=C)C=C5C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.4 (M + H)	3.84
3067	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=C(C=C4)C(=C)C=C5C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	3.46
3068	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=C(C=C4)C(=C)C=C5C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.4 (M + H)	3.79
3069	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=C(C=C4)C(=C)C=C5C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	3.55
3070	 <chem>CC1=NC2=CC=CC=C2N(C)=NC1NCC3CCCCC3NC(=O)C4=CC=C(C=C4)C5=C(C=C4)C(=C)C=C5C6=CC=CC=C6</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	3.63

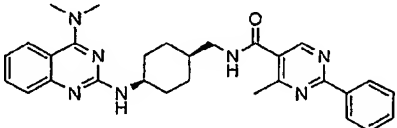
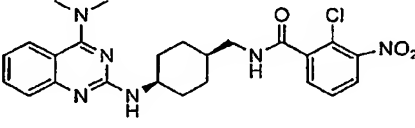
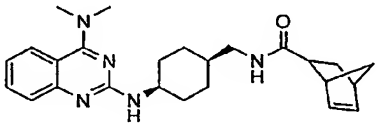
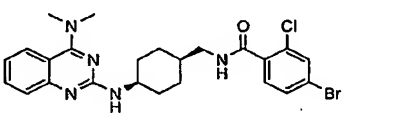
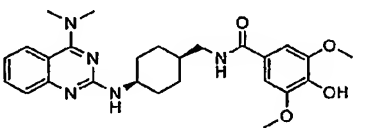
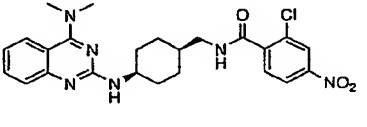
Example No.	Structure	ESI-MS	Retention Time (min)
3071	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4cc(F)c(F)c(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.69
3072	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4cc(F)c(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	3.69
3073	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4cc(F)c(F)c(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	3.66
3074	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4ccccc4F</chem> $\text{CF}_3\text{CO}_2\text{H}$	422.4 (M + H)	3.55
3075	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4ccc(CCC)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	4.24
3076	 <chem>CC1=NC2=CC=CC=C2N1C(C)N[C@H]3CCCC[C@H]3CNC(=O)c4ccc(C#N)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	429.2 (M + H)	3.42

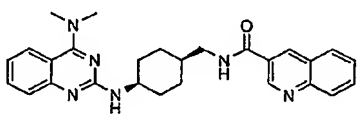
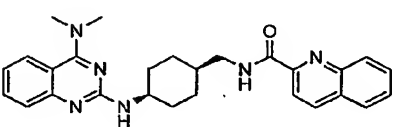
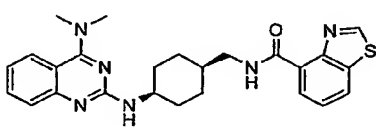
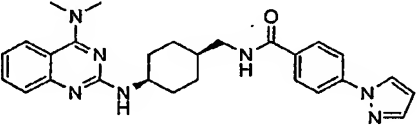
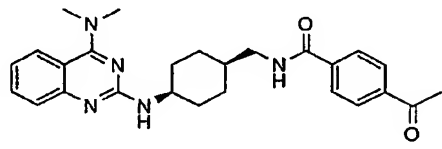
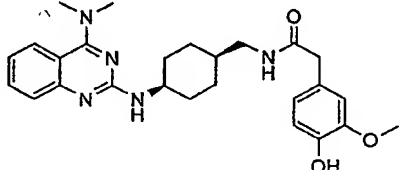
Example No.	Structure	ESI-MS	Retention Time (min)
3077	 <chem>COc1ccc(cc1)C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3</chem> $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.61
3078	 <chem>COc1ccc(cc1)C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	3.86
3079	 <chem>CCCCCOc1ccc(cc1)C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.74
3080	 <chem>Fc1cc(F)c(F)cc1C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.68
3081	 <chem>C1CCCCC1C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3</chem> $\text{CF}_3\text{CO}_2\text{H}$	410.4 (M + H)	3.58
3082	 <chem>Cc1cc(C)c(cc1C(=O)N[C@H]2CCCC[C@H]2Nc3nc(C)c(C)n3)C(F)(F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	540.4 (M + H)	4.19

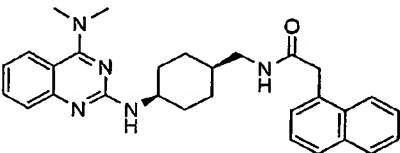
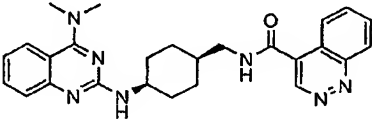
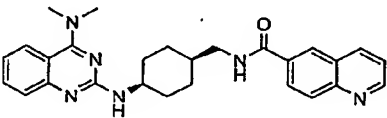
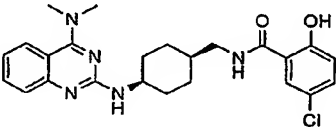
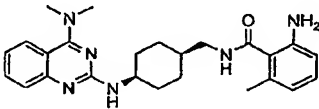
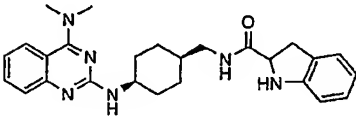
Example No.	Structure	ESI-MS	Retention Time (min)
3083	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3CNC(=O)c4ccc(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	422.2 (M + H)	3.50
3084	 <chem>COc1cc(OC)c(OC)c(C(=O)N[C@H]2CCCC[C@H]2CNC3=NC4=CC=CC=C4N(C)C3=NC4)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.4 (M + H)	3.39
3085	 <chem>FC1=CC=C(C(=C1)C(=O)N[C@H]2CCCC[C@H]2CNC3=NC4=CC=CC=C4N(C)C3=NC4)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.0 (M + H)	3.55
3086	 <chem>Clc1ccccc1C(=O)N[C@H]2CCCC[C@H]2CNC3=NC4=CC=CC=C4N(C)C3=NC4</chem> $\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	3.48
3087	 <chem>c1ccc2cc(C(=O)N[C@H]3CCCC[C@H]3CNC4=NC5=CC=CC=C5N(C)C4=NC5)c3ccccc23</chem> $\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	3.75
3088	 <chem>Cc1ccc(C(=O)N[C@H]2CCCC[C@H]2CNC3=NC4=CC=CC=C4N(C)C3=NC4)cc1C(F)(F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	3.83

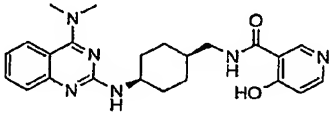
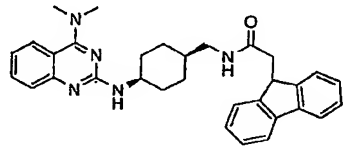
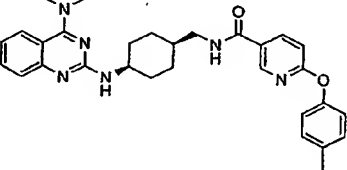
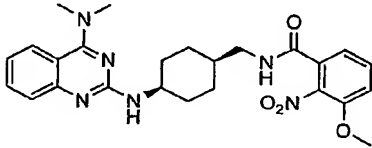
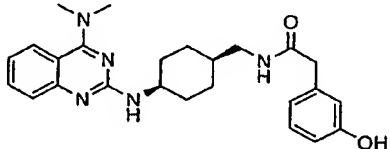
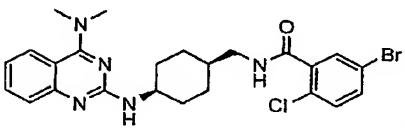
Example No.	Structure	ESI-MS	Retention Time (min)
3089	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC=C(C=C4)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	422.2 (M + H)	3.51
3090	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC=C(C=C4)C(F)(F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	3.87
3091	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC(=C(C=C4)F)Br</chem> $\text{CF}_3\text{CO}_2\text{H}$	500.4 (M + H)	3.03
3092	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC=C(C=C4)N(C)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	447.4 (M + H)	2.59
3093	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC=C(C=C4)c5ccsc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.4 (M + H)	3.25
3094	 <chem>CC1=NC2=CC=CC=C2N1C(C)C3CCCCC3NC(=O)C4=CC=C(C=C4)c5nn[nH]s5</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	2.81

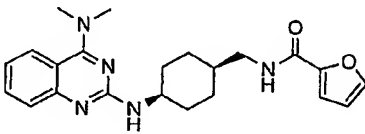
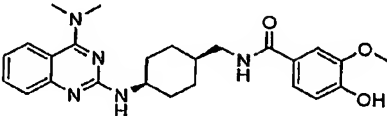
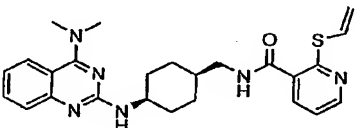
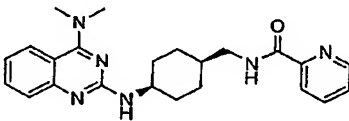
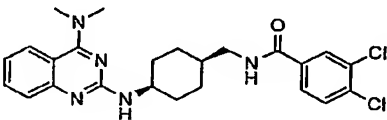
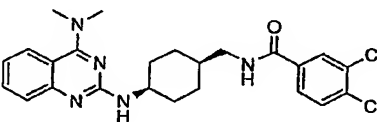
Example No.	Structure	ESI-MS	Retention Time (min)
3095	 <chem>CC1=CC=C(C=C1C(=O)NCC2CCCCC2N3C=NC4=CC=CC=C4N(C)C3=C)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	452.4 (M + H)	2.98
3096	 <chem>CC1=CC=C(C=C1C(=O)NCC2CCCCC2N3C=NC4=CC=CC=C4N(C)C3=C)COC5=CC=CC=C5</chem> $\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	3.29
3097	 <chem>CC(=O)NCC1CCCCC1N2C=NC3=CC=CC=C3N(C)C2=C1C</chem> $\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	2.77
3098	 <chem>O=C1C2=CC=CC=C2CCC1NCC3CCCCC3N4C=NC5=CC=CC=C5N(C)C4=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.4 (M + H)	3.06
3099	 <chem>CC1=CC=C(C=C1C(=O)NCC2CCCCC2N3C=NC4=CC=CC=C4N(C)C3=C)C2=CN=C(C2)C3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	484.4 (M + H)	3.40
3100	 <chem>CC(=O)NCC1CCCCC1N2C=NC3=CC=CC=C3N(C)C2=C1</chem> $\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	2.69

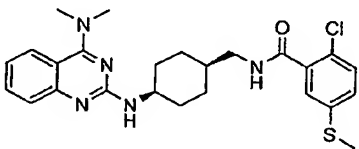
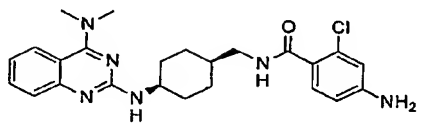
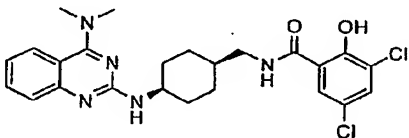
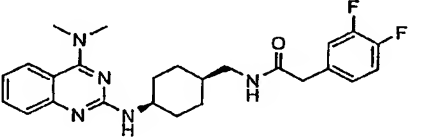
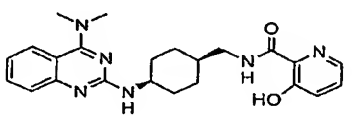
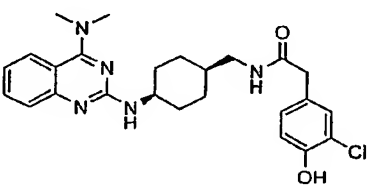
Example No.	Structure	ESI-MS	Retention Time (min)
3101	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)Nc4nc(C)c(C5=CC=CC=C5)n4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	3.01
3102	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)Nc4cc([N+](=O)[O-])cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	483.4 (M + H)	2.79
3103	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)N4C5C6C7C8C5C6C7C8C4</chem> $\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.76
3104	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)Nc4cc(Br)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.2 (M + H)	3.03
3105	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)Nc4cc(OC)c(O)c(OC)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	2.41
3106	 <chem>CC1=CN(C)C(=N1)N[C@H]2CCCC[C@H]2CN[C@@H]3CCCC[C@H]3C(=O)Nc4cc([N+](=O)[O-])cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	483.2 (M + H)	2.84

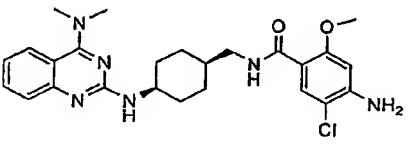
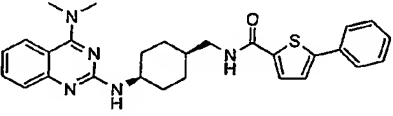
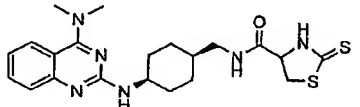
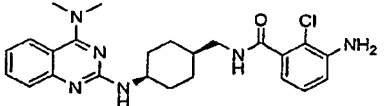
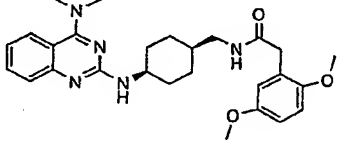
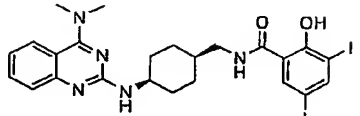
Example No.	Structure	ESI-MS	Retention Time (min)
3107	 $2\text{CF}_3\text{CO}_2\text{H}$	455 (M + H)	2.45
3108	 $2\text{CF}_3\text{CO}_2\text{H}$	455.2 (M + H)	3.19
3109	 $\text{CF}_3\text{CO}_2\text{H}$	461.4 (M + H)	2.60
3110	 $2\text{CF}_3\text{CO}_2\text{H}$	470.4 (M + H)	2.74
3111	 $\text{CF}_3\text{CO}_2\text{H}$	446.6 (M + H)	2.61
3112	 $\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	2.35

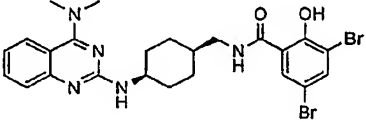
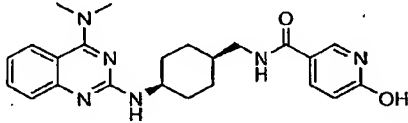
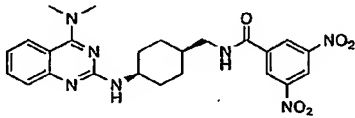
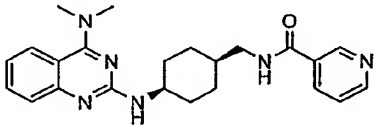
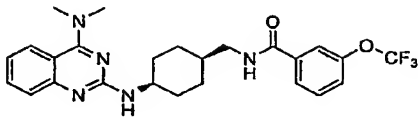
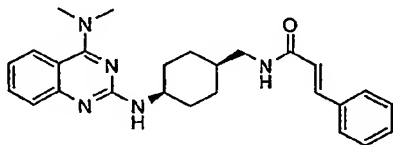
Example No.	Structure	ESI-MS	Retention Time (min)
3113	 <p>CF₃CO₂H</p>	468.4 (M + H)	3.04
3114	 <p>2CF₃CO₂H</p>	456.2 (M + H)	2.44
3115	 <p>2CF₃CO₂H</p>	455.2 (M + H)	2.11
3116	 <p>CF₃CO₂H</p>	454.2 (M + H)	3.21
3117	 <p>2CF₃CO₂H</p>	433.6 (M + H)	2.34
3118	 <p>2CF₃CO₂H</p>	444.6 (M+)	2.93

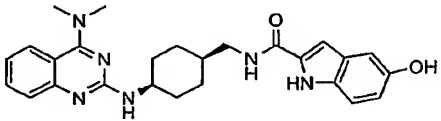
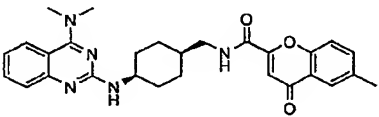
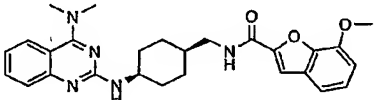
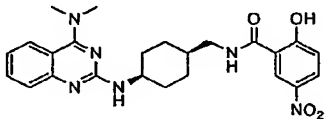
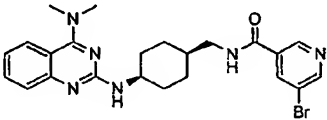
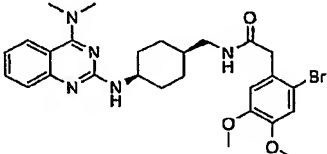
Example No.	Structure	ESI-MS	Retention Time (min)
3119	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)c4ccncc4C(=O)O</chem> $2\text{CF}_3\text{CO}_2\text{H}$	421.4 (M + H)	2.23
3120	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)Cc4c5ccccc5c6ccccc46C(=O)O</chem> $\text{CF}_3\text{CO}_2\text{H}$	506.4 (M + H)	3.31
3121	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)Cc4ccncc4Oc5ccc(OC)cc5C(=O)O</chem> $2\text{CF}_3\text{CO}_2\text{H}$	511.6 (M + H)	3.21
3122	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)Cc4ccncc4C(=O)c5cc(OC)cc([N+](=O)[O-])cc5C(=O)O</chem> $\text{CF}_3\text{CO}_2\text{H}$	479.4 (M + H)	3.60
3123	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)Cc4ccncc4C(=O)Cc5ccc(O)cc5C(=O)O</chem> $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	2.37
3124	 <chem>CN1C=NC2=C(N1)C(=O)N(C2)C3CCCCC3NC(=O)Cc4ccncc4C(=O)c5cc(Br)cc(Cl)cc5C(=O)O</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.4 (M + H)	3.02

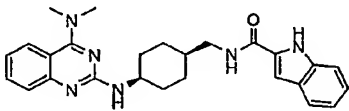
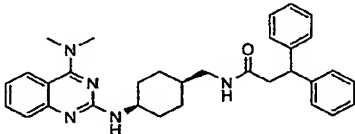
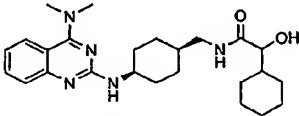
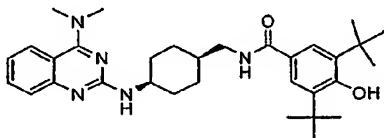
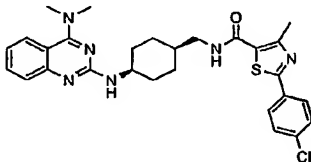
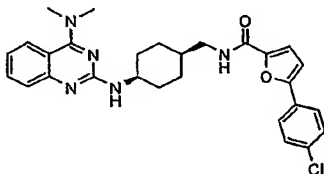
Example No.	Structure	ESI-MS	Retention Time (min)
3125	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4ccoc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	394.4 (M + H)	2.45
3126	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4cc(OC)c(O)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.2 (M + H)	2.41
3127	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4cc(C=C)sc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	477.0 (M + H)	2.88
3128	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4ccccn4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	405.6 (M + H)	2.61
3129	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4cc(Cl)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.6 (M + H)	3.17
3130	 <chem>CC1=NC2=CC=CC=C2N1C(C)=N[C@H]3CCCC[C@H]3CNC(=O)c4cc(Cl)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	2.59

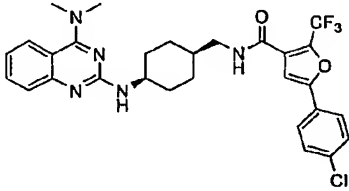
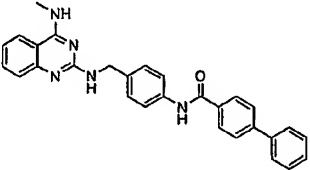
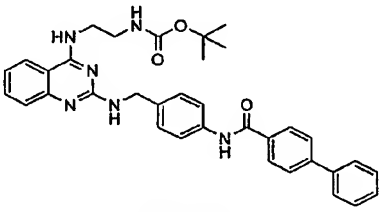
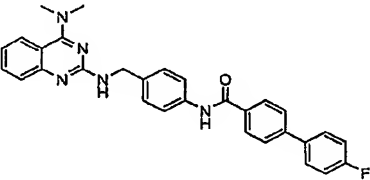
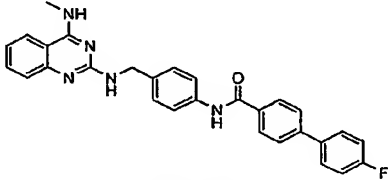
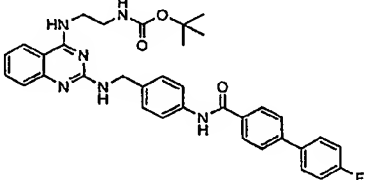
Example No.	Structure	ESI-MS	Retention Time (min)
3131	 $\text{CF}_3\text{CO}_2\text{H}$	484.2 (M + H)	2.99
3132	 $2\text{CF}_3\text{CO}_2\text{H}$	453.0 (M + H)	2.45
3133	 $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	3.59
3134	 $\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	2.81
3135	 $2\text{CF}_3\text{CO}_2\text{H}$	421.4 (M + H)	2.89
3136	 $\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	2.53

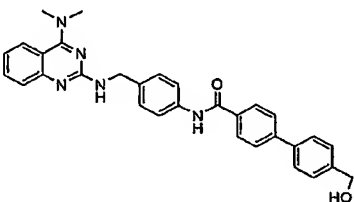
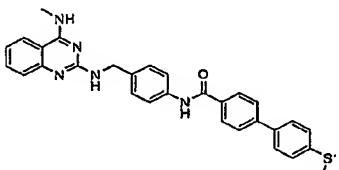
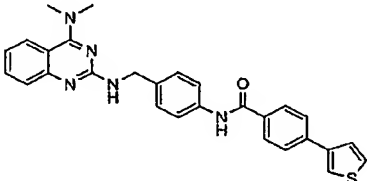
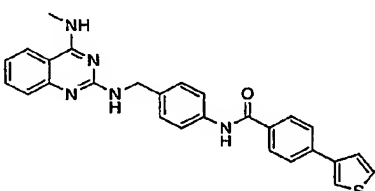
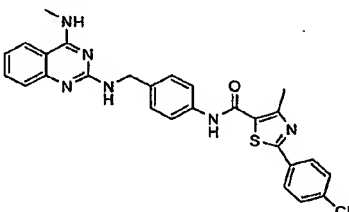
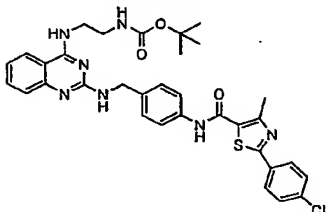
Example No.	Structure	ESI-MS	Retention Time (min)
3137	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)c2cc(N)c(Cl)c(OC)c2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	483.2 (M + H)	2.83
3138	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)c1cc2ccccc2s1</chem> $\text{CF}_3\text{CO}_2\text{H}$	487.4 (M+2H+)	3.40
3139	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)C1=NC(=S)SC1</chem> $\text{CF}_3\text{CO}_2\text{H}$	445.6 (M + H)	2.36
3140	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)c1cc(N)c(Cl)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	453.2 (M + H)	2.46
3141	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)Cc1cc(OC)c(OC)c(OC)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	2.77
3142	 <chem>CC1=NC2=CC=CC=C2N1N(C)CNC1CCC(CC1)CNC(=O)c1cc(O)c(I)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	672.2 (M + H)	3.92

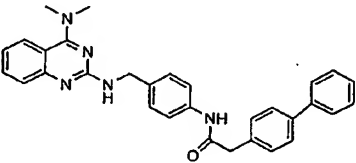
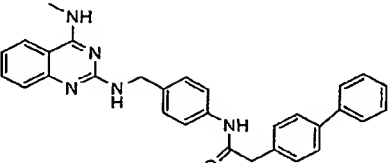
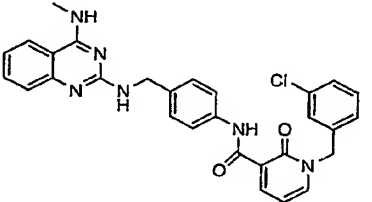
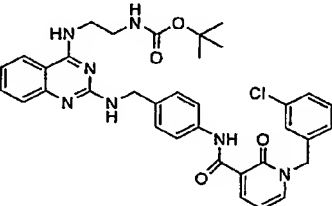
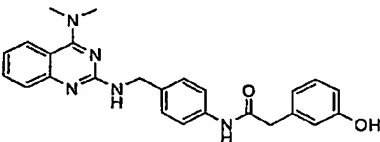
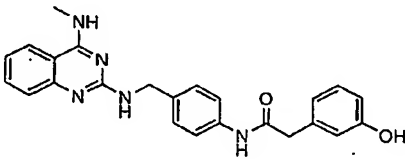
Example No.	Structure	ESI-MS	Retention Time (min)
3143	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)c1cc(Br)cc(Br)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	576.2 (M + H)	3.71
3144	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)c1cccc(O)c1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	421.2 (M + H)	2.01
3145	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)c1cc([N+](=O)[O-])cc([N+](=O)[O-])c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.4 (M + H)	2.77
3146	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)c1ccccn1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	405.6 (M + H)	1.99
3147	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)c1ccc(OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	3.13
3148	 <chem>CC1=NC2=CC=CC=C2N1N[C@H]3CCCC[C@H]3CNC(=O)/C=C/c1ccccc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	430.4 (M + H)	2.91

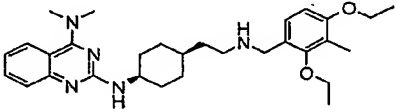
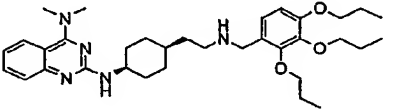
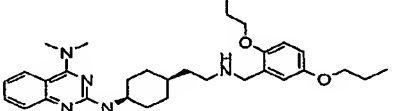
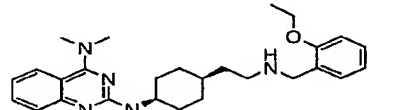
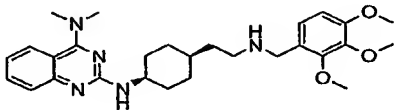
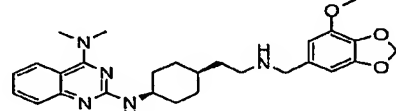
Example No.	Structure	ESI-MS	Retention Time (min)
3149	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4c[nH]c5ccc(O)cc45)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	459.4 (M + H)	2.47
3150	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4c5cc(=O)ccc5oc4=O)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.6 (M + H)	2.93
3151	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4c5cc(OC)ccc5oc4=O)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	3.03
3152	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4ccc([N+](=O)[O-])cc4)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	465.2 (M + H)	3.13
3153	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4ccncc4Br)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	483.4 (M + H)	2.67
3154	 <chem>CC1=CN(C)N=C2C=CC=CC=C12N[C@H]3CCCC[C@H]3CN(C(=O)Nc4cc(OC)c(OC)c(Br)c4)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	556.4 (M + H)	2.84

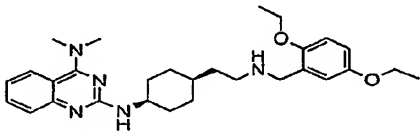
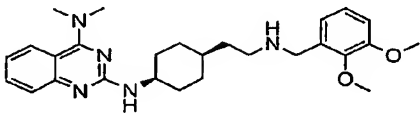
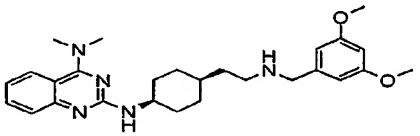
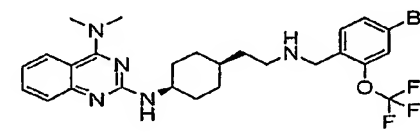
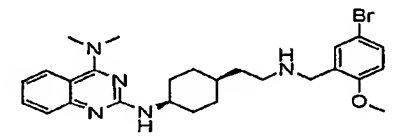
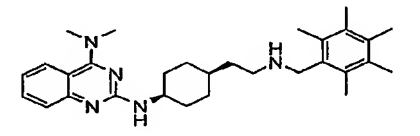
Example No.	Structure	ESI-MS	Retention Time (min)
3155	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)Nc4ccccc4</chem> <chem>2CF3CO2H</chem>	443.4 (M + H)	2.94
3156	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)N(Cc4ccccc4)Cc5ccccc5</chem> <chem>CF3CO2H</chem>	508.2 (M + H)	3.20
3157	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)NCC4CCCCC4C(=O)O</chem> <chem>CF3CO2H</chem>	440.0 (M + H)	2.72
3158	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)NCC4C(C)(C)C(=O)C(C)(C)C4O</chem> <chem>CF3CO2H</chem>	532.4 (M + H)	3.58
3159	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)Nc4nc5cc(Cl)ccc5s4</chem> <chem>CF3CO2H</chem>	535.4 (M + H)	3.51
3160	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)Nc4oc5cc(Cl)ccc5o4</chem> <chem>CF3CO2H</chem>	504.4 (M + H)	3.49

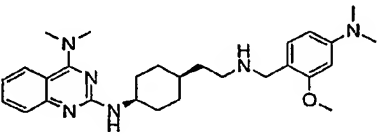
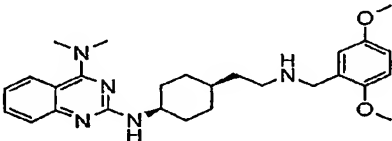
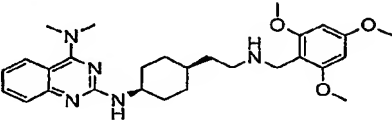
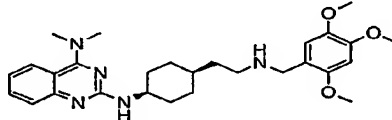
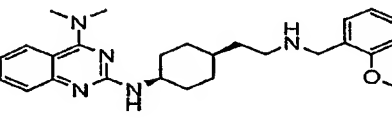
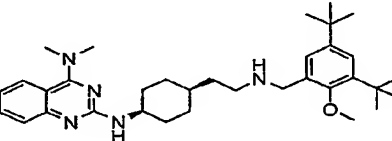
Example No.	Structure	ESI-MS	Retention Time (min)
3161	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3CCCCC3CC(=O)c4oc(cc4C(F)(F)F)c5ccc(Cl)cc5</chem> <chem>CF3CO2H</chem>	572.4 (M + H)	3.71
3162	 <chem>CC1=NC2=CC=CC=C2N=C1NCC3=CC=CC=C3CC(=O)Nc4ccc(cc4)-c5ccccc5</chem> <chem>CF3CO2H</chem>	460.2 (M + H)	3.80
3163	 <chem>CC(C)(C)OC(=O)NCC1=NC2=CC=CC=C2N=C1NCC3=CC=CC=C3CC(=O)Nc4ccc(cc4)-c5ccccc5</chem> <chem>CF3CO2H</chem>	589.2 (M + H)	4.00
3164	 <chem>CC1=NC2=CC=CC=C2N(C)N=C1NCC3=CC=CC=C3CC(=O)Nc4ccc(cc4)-c5ccc(F)cc5</chem> <chem>CF3CO2H</chem>	492.2 (M + H)	3.90
3165	 <chem>CC1=NC2=CC=CC=C2N=C1NCC3=CC=CC=C3CC(=O)Nc4ccc(cc4)-c5ccc(F)cc5</chem> <chem>CF3CO2H</chem>	478.2 (M + H)	3.80
3166	 <chem>CC(C)(C)OC(=O)NCC1=NC2=CC=CC=C2N=C1NCC3=CC=CC=C3CC(=O)Nc4ccc(cc4)-c5ccc(F)cc5</chem> <chem>CF3CO2H</chem>	607.6 (M + H)	4.00

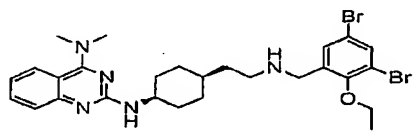
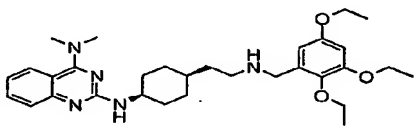
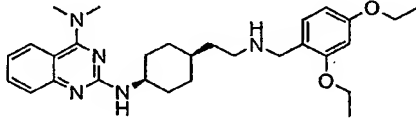
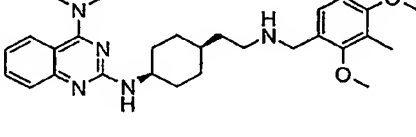
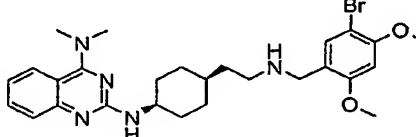
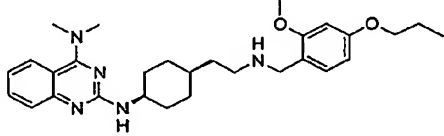
Example No.	Structure	ESI-MS	Retention Time (min)
3167	 <chem>CC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC=C(CO)C=C4)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	504.2 (M + H)	3.40
3168	 <chem>CC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC=C(CI)C=C4)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	506.2 (M + H)	3.90
3169	 <chem>CC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC=C(C5=CC=CS5)C=C4)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	3.80
3170	 <chem>CC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC(=C(C5=CC=CS5))C=C4)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	466.2 (M + H)	3.70
3171	 <chem>CC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC(=CC=C4)C5=NC6=CC(=CC=C6)S=C5N6C7=CC=CC=C7Cl)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	515.2 (M + H)	3.90
3172	 <chem>CC(C)(C)OC(=O)NCC1=NC2=CC=CC=C2N1N=C(NCC3=CC=C(NC(=O)C4=CC(=CC=C4)C5=NC6=CC(=CC=C6)S=C5N6C7=CC=CC=C7Cl)C=C3)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	644.2 (M + H)	4.10

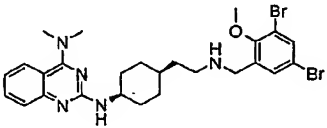
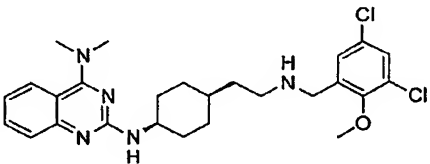
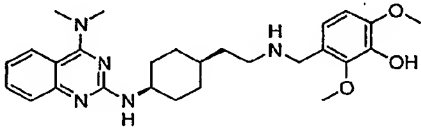
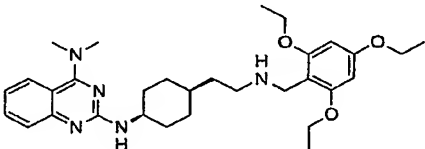
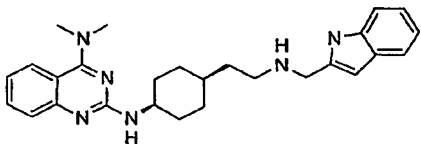
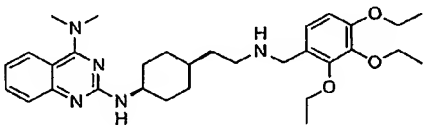
Example No.	Structure	ESI-MS	Retention Time (min)
3173	 <chem>CC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.2 (M + H)	3.90
3174	 <chem>CC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=CC=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	3.80
3175	 <chem>CC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=C(Cl)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	525.4 (M + H)	3.70
3176	 <chem>CC(C)(C)OC(=O)NCC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=C(Cl)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	654.2 (M + H)	3.90
3177	 <chem>CC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=C(O)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	428.2 (M + H)	3.10
3178	 <chem>CC1=NC2=CC=CC=C2N1NCC(=O)NCC3=CC=C(O)C=C3</chem> $\text{CF}_3\text{CO}_2\text{H}$	414.4 (M + H)	2.90

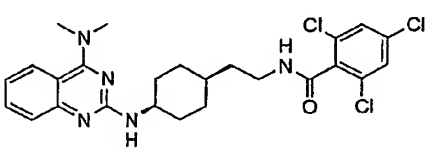
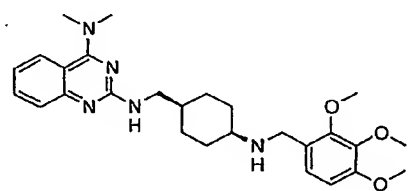
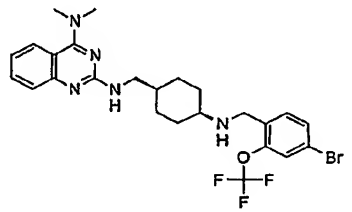
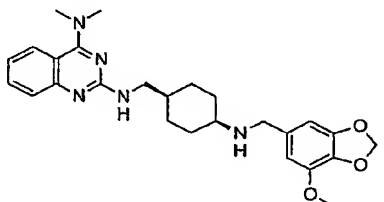
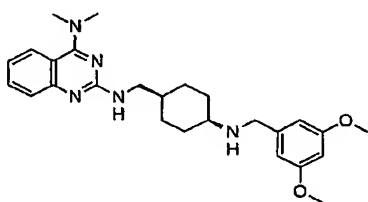
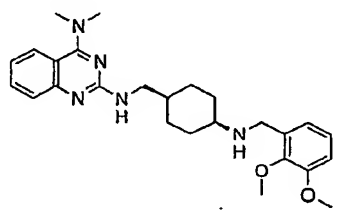
Example No.	Structure	ESI-MS	Retention Time (min)
3179	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	506.4 (M + H)	3.04
3180	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	578.8 (M + H)	3.50
3181	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	520.6 (M + H)	3.19
3182	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	448.4 (M + H)	2.80
3183	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	494.6 (M + H)	2.66
3184	 <chem>CCOC1=CC=C(C(=C1)NCC2=CC=CC=C2N2C(=N3C(=N2)C(=C4C=CC=CC=C4N(C)C)N3)N2)CC2</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	2.66

Example No.	Structure	ESI-MS	Retention Time (min)
3185	 <chem>CCOC1=CC=C(C(=C1)OCC)NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.6 (M + H)	2.94
3186	 <chem>COc1ccc(NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5)cc1OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	2.65
3187	 <chem>COc1ccc(NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5)cc1OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	2.68
3188	 <chem>Brc1ccc(NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5)cc1OC(F)(F)F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	566.4 (M + H)	3.03
3189	 <chem>Brc1ccc(NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5)cc1OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	512.6 (M + H)	2.85
3190	 <chem>Cc1c(C)ccc(NCCN2CCCCC2C3=NC4=CC=CC=C4N(C)C3=NC5=CC=CC=C5N5)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	3.09

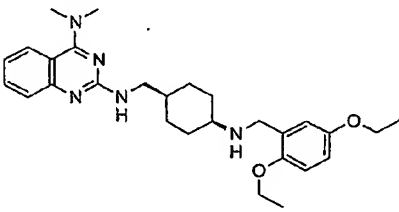
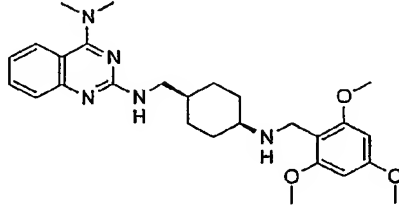
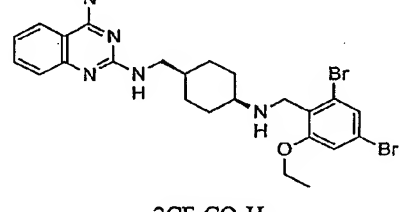
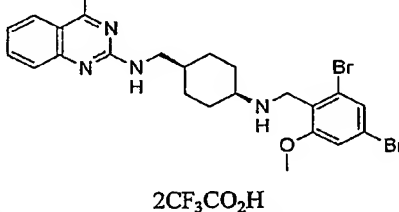
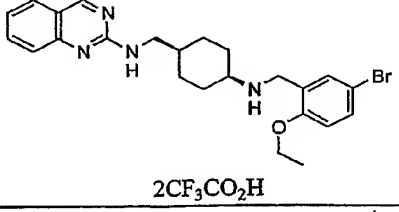
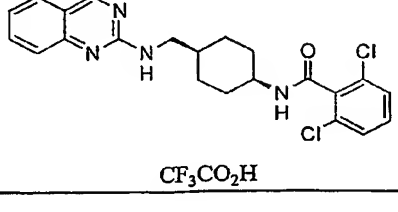
Example No.	Structure	ESI-MS	Retention Time (min)
3191	 $3\text{CF}_3\text{CO}_2\text{H}$	477.4 (M + H)	2.51
3192	 $2\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	2.67
3193	 $2\text{CF}_3\text{CO}_2\text{H}$	494.6 (M + H)	2.78
3194	 $2\text{CF}_3\text{CO}_2\text{H}$	494.6 (M + H)	2.60
3195	 $2\text{CF}_3\text{CO}_2\text{H}$	434.6 (M + H)	2.67
3196	 $2\text{CF}_3\text{CO}_2\text{H}$	546.4 (M + H)	4.30

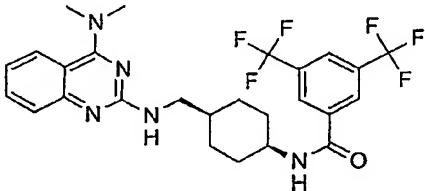
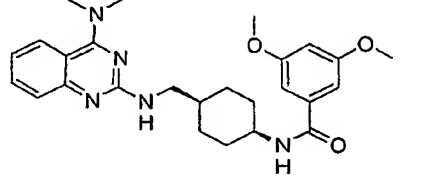
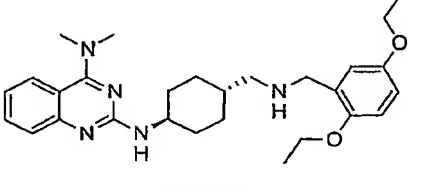
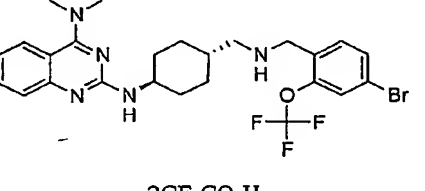
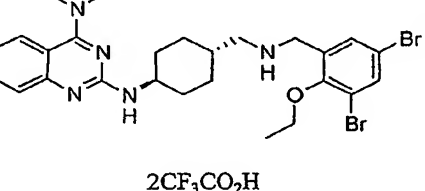
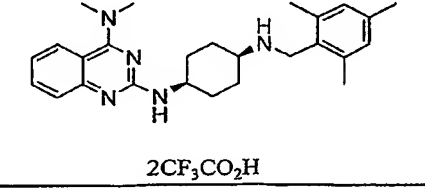
Example No.	Structure	ESI-MS	Retention Time (min)
3197	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4cc(Br)c(Br)oc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	606.6 (M + H)	3.95
3198	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4cc(OC)c(OC)c(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	536.6 (M + H)	3.83
3199	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4cc(OC)c(OCC)c(OCC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.4 (M + H)	2.97
3200	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4cc(OC)c(C)c(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	2.79
3201	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4cc(Br)c(OC)c(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	542.0 (M + H)	2.85
3202	 <chem>CC1=CC=C2N=CN(C1)N2C3CCCCC3CNCCc4ccc(OC)c(OCC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	492.6 (M + H)	2.81

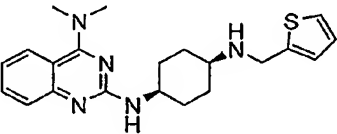
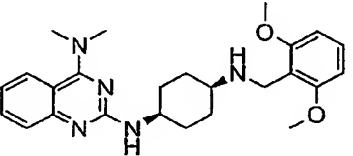
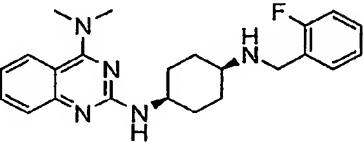
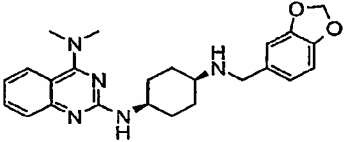
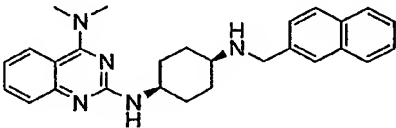
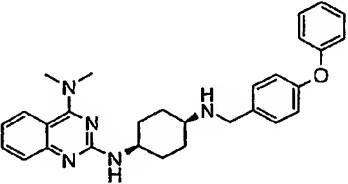
Example No.	Structure	ESI-MS	Retention Time (min)
3203	 <chem>COc1cc(Br)cc(Br)cc1NCCN2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	590.4 (M + H)	3.02
3204	 <chem>COc1cc(Cl)cc(Cl)cc1NCCN2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	502.2 (M + H)	2.91
3205	 <chem>COc1cc(OC)cc(O)cc1NCCN2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	2.51
3206	 <chem>CCOc1cc(OC)cc(OC)cc1NCCN2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	3.21
3207	 <chem>c1ccc2c(c1)c(c[nH]2)CNCCN3CCCCC3Nc4nc5ccccc5n4C</chem> $3\text{CF}_3\text{CO}_2\text{H}$	443.6 (M + H)	2.66
3208	 <chem>CCOc1cc(OC)cc(OC)cc1NCCN2CCCCC2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	3.08

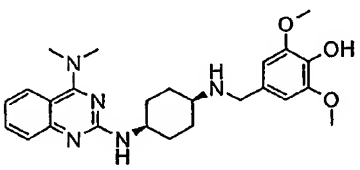
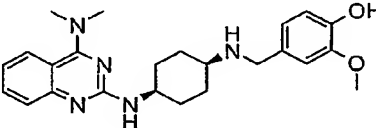
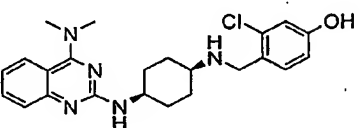
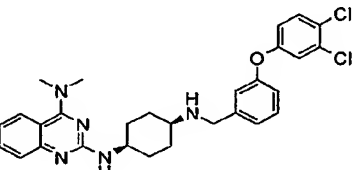
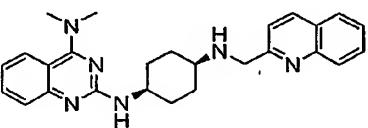
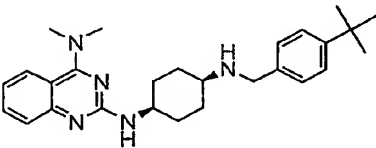
Example No.	Structure	ESI-MS	Retention Time (min)
3209	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC=C(C(=O)N)C(Cl)=CC5Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	520.0 (M + H)	3.51
3210	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC(OC)=C(OC)C(OC)=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	2.58
3211	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC=C(C(F)(F)F)C(Br)=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	552.0 (M + H)	3.11
3212	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC6=C(C=C5)OCOC6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	3.22
3213	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC(OC)=CC(OC)=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	2.70
3214	 <chem>CC1=NC2=C(N1)N=CN=C2C3=CC=CC=C3NCC4CCCCC4NCC5=CC(OC)=CC(OC)=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	2.58

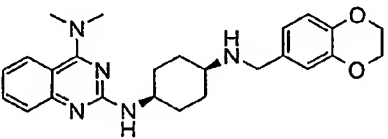
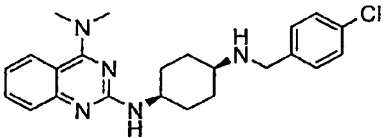
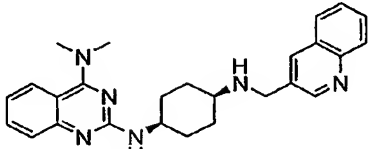
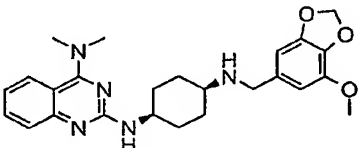
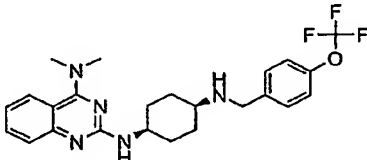
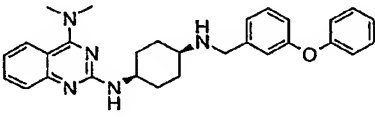
Example No.	Structure	ESI-MS	Retention Time (min)
3215	 <chem>COc1ccc(cc1)CNCC2CCCCC2CNCC3=NC(=N)N(C)C3c4ccccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	2.73
3216	 <chem>c1ccc2c(c1)c[nH]2CNCC3CCCCC3CNCC4=NC(=N)N(C)C4c5ccccc5</chem> $3\text{CF}_3\text{CO}_2\text{H}$	429.4 (M + H)	3.29
3217	 <chem>COc1cc(OC)c(OC)cc1CNCC2CCCCC2CNCC3=NC(=N)N(C)C3c4ccccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	2.78
3218	 <chem>CCOc1cc(OC)cc(OC)c1CNCC2CCCCC2CNCC3=NC(=N)N(C)C3c4ccccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	522.4 (M + H)	3.77
3219	 <chem>COc1ccccc1CNCC2CCCCC2CNCC3=NC(=N)N(C)C3c4ccccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	450.2 (M + H)	2.57
3220	 <chem>COc1ccc(Br)cc1CNCC2CCCCC2CNCC3=NC(=N)N(C)C3c4ccccc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	498.0 (M + H)	2.97

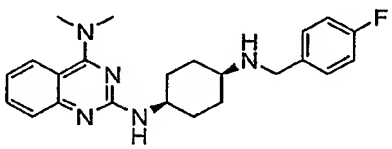
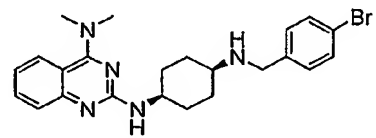
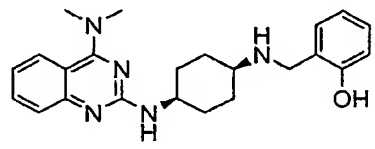
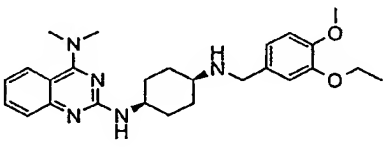
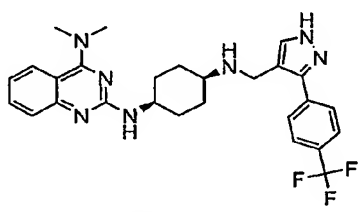
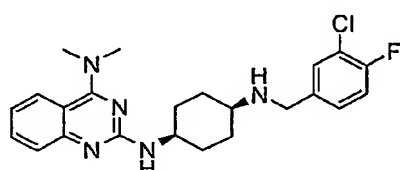
Example No.	Structure	ESI-MS	Retention Time (min)
3221	 $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	3.17
3222	 $2\text{CF}_3\text{CO}_2\text{H}$	480.0 (M + H)	3.08
3223	 $2\text{CF}_3\text{CO}_2\text{H}$	590.2 (M + H)	4.20
3224	 $2\text{CF}_3\text{CO}_2\text{H}$	576.4 (M + H)	3.95
3225	 $2\text{CF}_3\text{CO}_2\text{H}$	512.4 (M + H)	3.86
3226	 $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	3.07

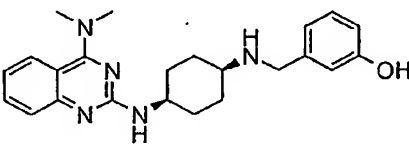
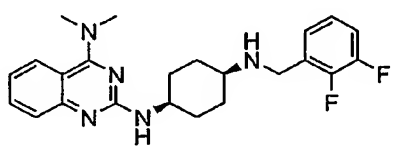
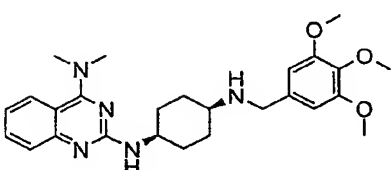
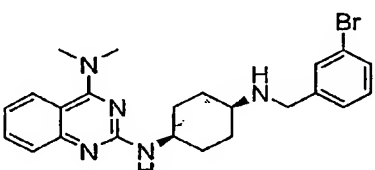
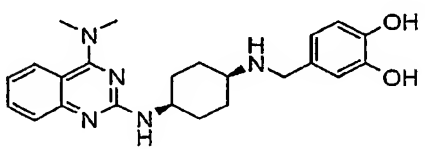
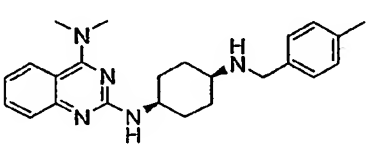
Example No.	Structure	ESI-MS	Retention Time (min)
3227	 $\text{CF}_3\text{CO}_2\text{H}$	540.6 (M + H)	3.75
3228	 $\text{CF}_3\text{CO}_2\text{H}$	464.4 (M + H)	3.07
3229	 $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	3.40
3230	 $2\text{CF}_3\text{CO}_2\text{H}$	552.6 (M + H)	3.50
3231	 $2\text{CF}_3\text{CO}_2\text{H}$	590.2 (M + H)	3.60
3232	 $2\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	3.25

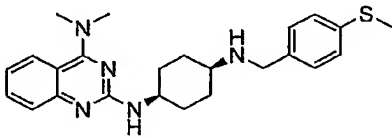
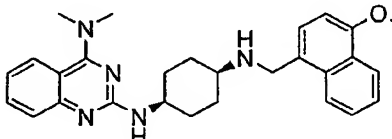
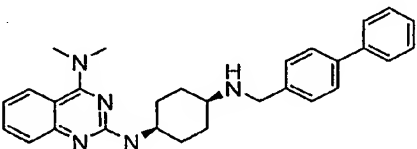
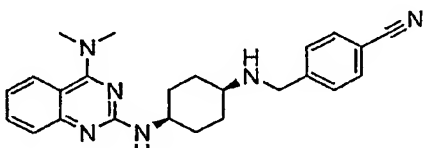
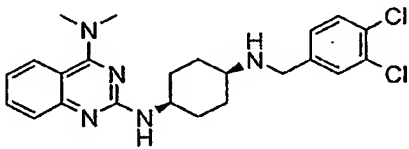
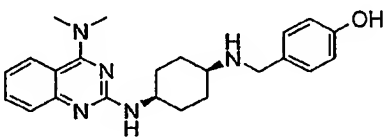
Example No.	Structure	ESI-MS	Retention Time (min)
3233	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4ccsc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	382.2 (M + H)	2.67
3234	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4cc(OC)c(OC)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	3.05
3235	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4ccccc4F</chem> $2\text{CF}_3\text{CO}_2\text{H}$	394.4 (M + H)	2.75
3236	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4cc5c(c(c4)OCO5)</chem> $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.82
3237	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4c5ccccc4c5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	426.4 (M + H)	3.17
3238	 <chem>CN(C)c1nc2c(ncn2C1)N[C@H]3CCCC[C@H]3NCc4ccc(Oc5ccccc5)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	3.44

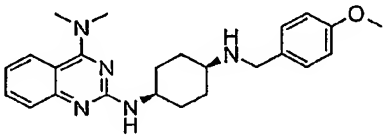
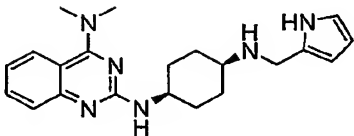
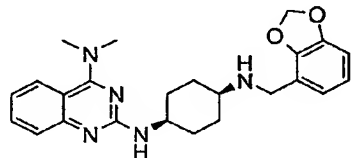
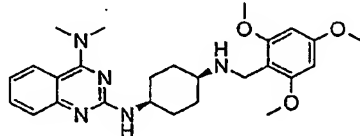
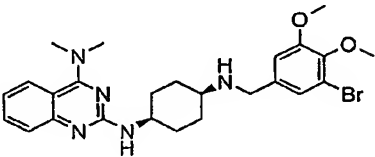
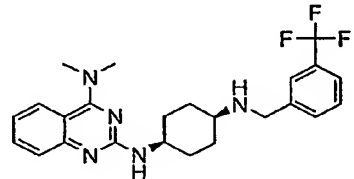
Example No.	Structure	ESI-MS	Retention Time (min)
3239	 $2\text{CF}_3\text{CO}_2\text{H}$	452.2 (M + H)	2.69
3240	 $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	2.80
3241	 $2\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	2.79
3242	 $2\text{CF}_3\text{CO}_2\text{H}$	536.4 (M + H)	3.75
3243	 $3\text{CF}_3\text{CO}_2\text{H}$	427.2 (M + H)	2.95
3244	 $2\text{CF}_3\text{CO}_2\text{H}$	432.4 (M + H)	3.41

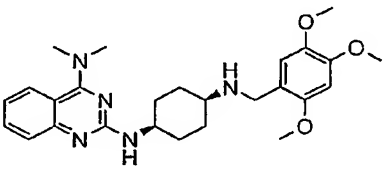
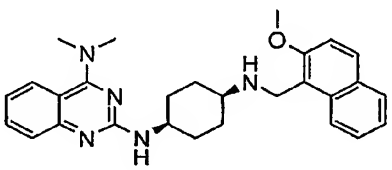
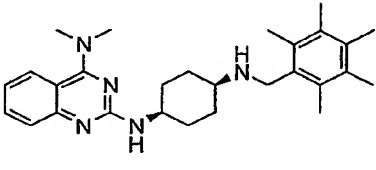
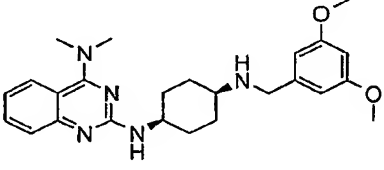
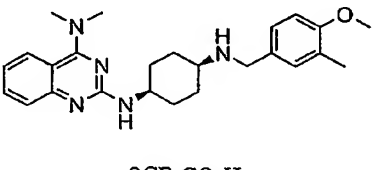

Example No.	Structure	ESI-MS	Retention Time (min)
3245	 $2\text{CF}_3\text{CO}_2\text{H}$	434.2 (M + H)	2.84
3246	 $2\text{CF}_3\text{CO}_2\text{H}$	410.2 (M + H)	3.02
3247	 $3\text{CF}_3\text{CO}_2\text{H}$	427.4 (M + H)	2.61
3248	 $2\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	2.91
3249	 $2\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	3.19
3250	 $2\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	2.79

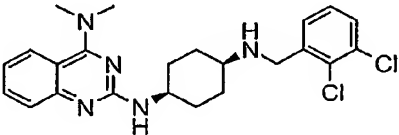
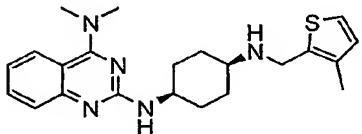
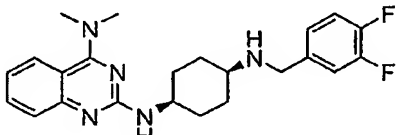
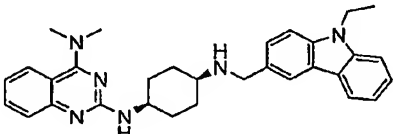
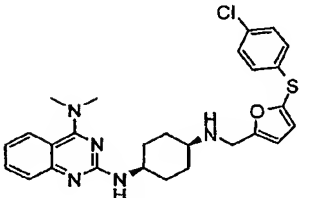
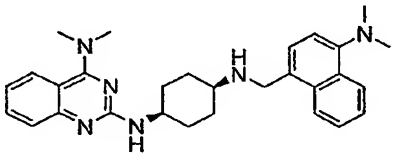
Example No.	Structure	ESI-MS	Retention Time (min)
3251	 $2\text{CF}_3\text{CO}_2\text{H}$	394.4 (M + H)	2.83
3252	 $2\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	3.08
3253	 $2\text{CF}_3\text{CO}_2\text{H}$	392.4 (M + H)	2.73
3254	 $2\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	2.92
3255	 $3\text{CF}_3\text{CO}_2\text{H}$	510.4 (M + H)	3.17
3256	 $2\text{CF}_3\text{CO}_2\text{H}$	428.2 (M + H)	3.08

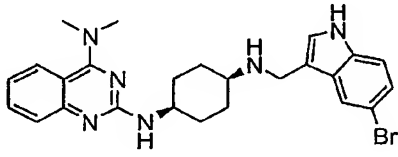
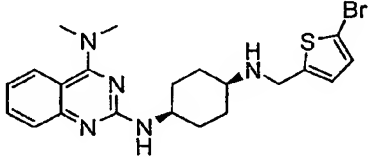
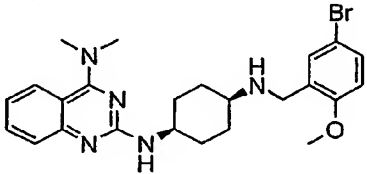
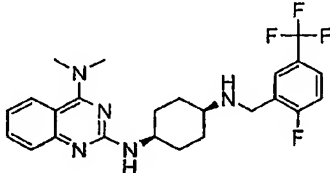
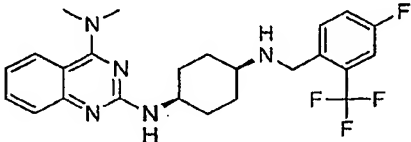
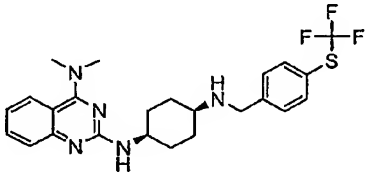
Example No.	Structure	ESI-MS	Retention Time (min)
3257	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4ccc(O)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	392.4 (M + H)	2.63
3258	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4cc(F)c(F)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	412.2 (M + H)	2.83
3259	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4cc(OC)c(OC)c(OC)c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	2.89
3260	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4ccc(Br)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	454.0 (M + H)	3.05
3261	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4cc(O)c(O)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	408.2 (M + H)	2.53
3262	 <chem>CN(C)c1nc2ccccc2n1N[C@H]3CCCC[C@H]3NCc4ccc(C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	390.4 (M + H)	2.92

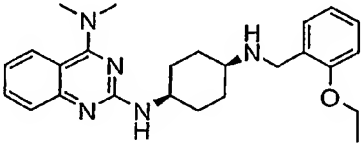
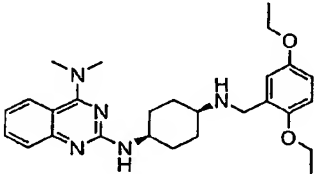
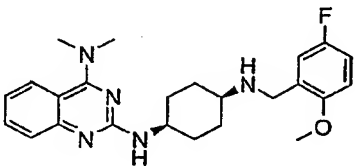
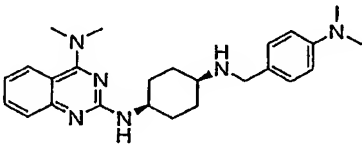
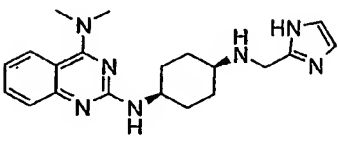
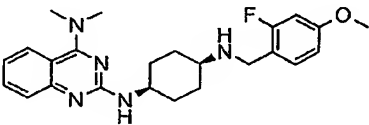
Example No.	Structure	ESI-MS	Retention Time (min)
3263	 <chem>CN(C)c1nc2c(ncn2C1NCCc3ccc(SC)cc3)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	422.2 (M + H)	3.05
3264	 <chem>CN(C)c1nc2c(ncn2C1NCCc3ccc(OC)cc3)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	456.4 (M + H)	3.25
3265	 <chem>CN(C)c1nc2c(ncn2C1NCCc3ccc(cc3)-c4ccccc4)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	452.2 (M + H)	3.37
3266	 <chem>CN(C)c1nc2c(ncn2C1NCCc3ccc(C#N)cc3)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	401.2 (M + H)	2.76
3267	 <chem>CN(C)c1nc2c(ncn2C1NCCc3cc(Cl)c(Cl)cc3)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	3.17
3268	 <chem>CN(C)c1nc2c(ncn2C1NCCc3ccc(O)cc3)c3ccccc3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	392.4 (M + H)	2.61

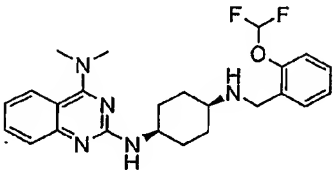
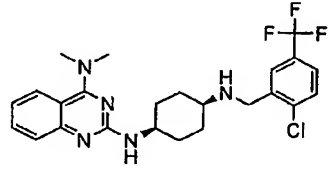
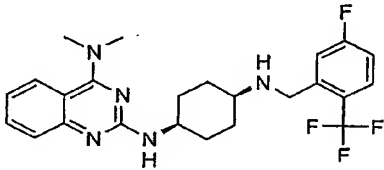
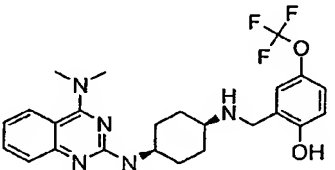
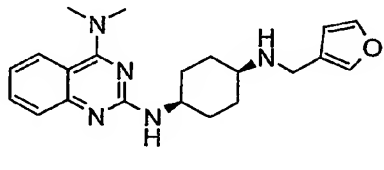
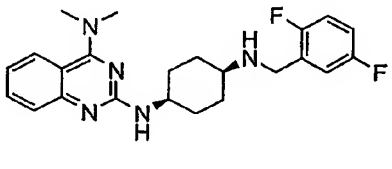
Example No.	Structure	ESI-MS	Retention Time (min)
3269	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4ccc(OC)cc4)c5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	406.4 (M + H)	2.86
3270	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4c[nH]c5ccccc45)c6ccccc16</chem> $3\text{CF}_3\text{CO}_2\text{H}$	365.4 (M + H)	2.61
3271	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4c5ccccc5oc4O)c6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.83
3272	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4cc(OC)c(OC)c(OC)c4)c5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	3.10
3273	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4cc(OC)c(Br)cc4)c5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	514.4 (M + H)	3.13
3274	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NCc4ccc(C(F)(F)F)cc4)c5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	3.17

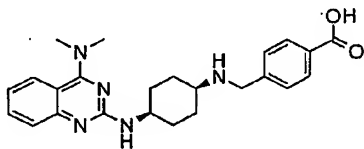
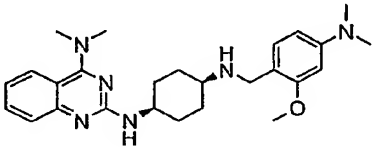
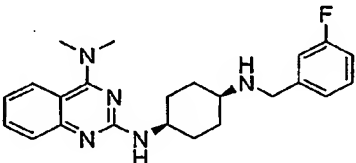
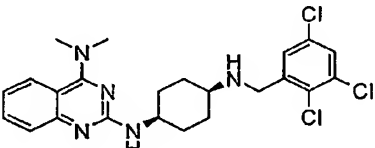
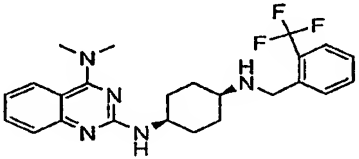
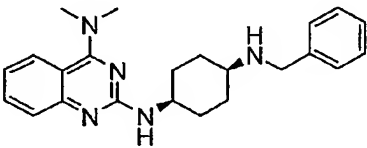
Example No.	Structure	ESI-MS	Retention Time (min)
3275	 2CF ₃ CO ₂ H	466.4 (M + H)	2.86
3276	 2CF ₃ CO ₂ H	456.2 (M + H)	3.22
3277	 2CF ₃ CO ₂ H	446.6 (M + H)	3.45
3278	 2CF ₃ CO ₂ H	436.4 (M + H)	2.95
3279	 2CF ₃ CO ₂ H	420.2 (M + H)	3.03
3280	 2CF ₃ CO ₂ H	382.4 (M + H)	2.72

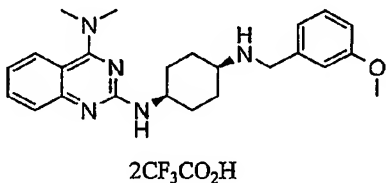
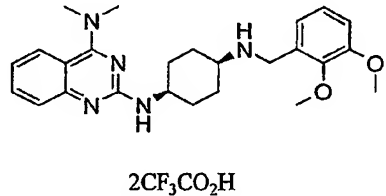
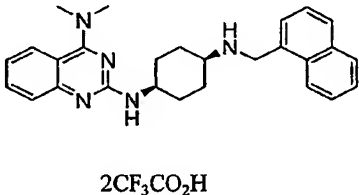
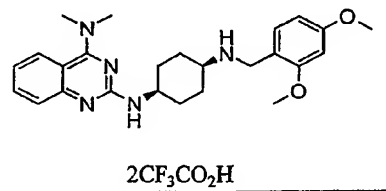
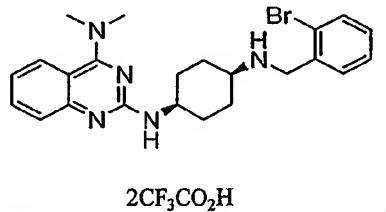
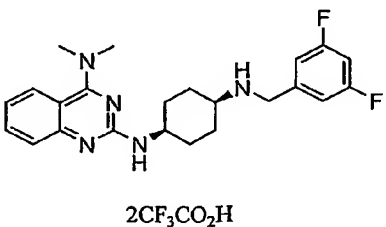
Example No.	Structure	ESI-MS	Retention Time (min)
3281	 $2\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	3.07
3282	 $2\text{CF}_3\text{CO}_2\text{H}$	396.2 (M + H)	2.79
3283	 $2\text{CF}_3\text{CO}_2\text{H}$	412.4 (M + H)	2.95
3284	 $32\text{CF}_3\text{CO}_2\text{H}$	493.4 (M + H)	3.57
3285	 $2\text{CF}_3\text{CO}_2\text{H}$	508.2 (M + H)	3.52
3286	 $2\text{CF}_3\text{CO}_2\text{H}$	469.6 (M + H)	2.76

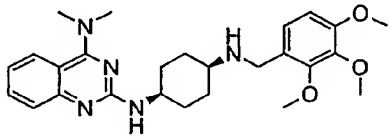
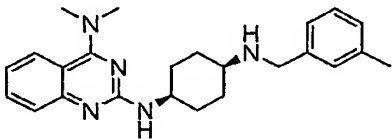
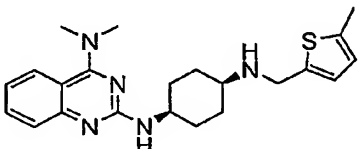
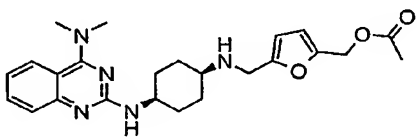
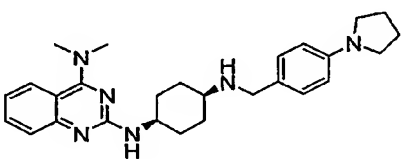
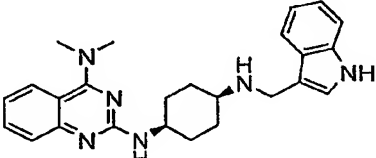
Example No.	Structure	ESI-MS	Retention Time (min)
3287	 $3\text{CF}_3\text{CO}_2\text{H}$	493.2 (M + H)	3.17
3288	 $2\text{CF}_3\text{CO}_2\text{H}$	460.2 (M + H)	2.95
3289	 $2\text{CF}_3\text{CO}_2\text{H}$	484.2 (M + H)	3.14
3290	 $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.11
3291	 $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.11
3292	 $2\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	3.39

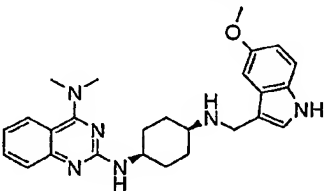
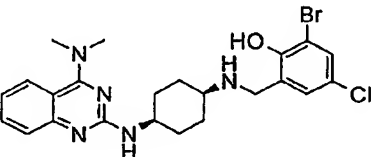
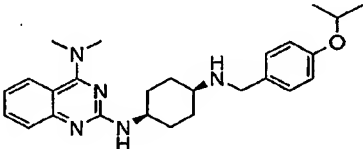
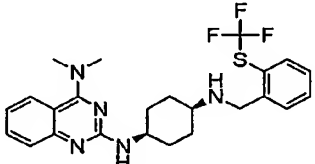
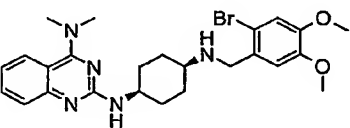
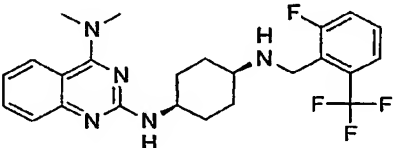
Example No.	Structure	ESI-MS	Retention Time (min)
3293	 $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	3.05
3294	 $2\text{CF}_3\text{CO}_2\text{H}$	464.2 (M + H)	3.21
3295	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	2.94
3296	 $3\text{CF}_3\text{CO}_2\text{H}$	419.4 (M + H)	2.51
3297	 $3\text{CF}_3\text{CO}_2\text{H}$	366.4 (M + H)	2.26
3298	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	2.93

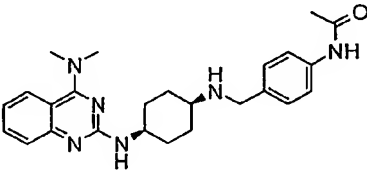
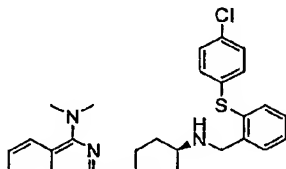
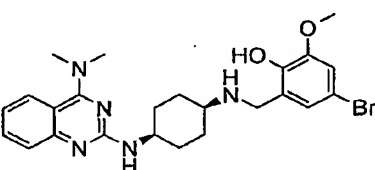
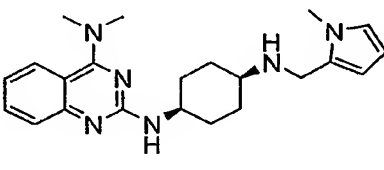
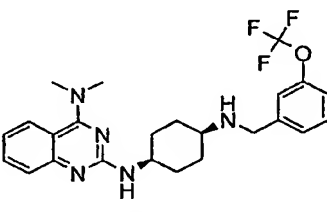
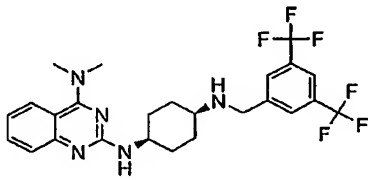
Example No.	Structure	ESI-MS	Retention Time (min)
3299	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccc(OC(F)(F)F)cc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	442.4 (M + H)	2.97
3300	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccc(C(F)(F)F)cc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.2 (M + H)	3.19
3301	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccc(F)cc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.05
3302	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccc(OC(F)(F)F)cc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	3.20
3303	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccoc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	366.4 (M + H)	2.64
3304	 <chem>CN(C)c1nc2c(nc3ccccc12)N[C@H]4CCCC[C@H]4N[C@@H](Cc1ccc(F)cc1)C5=CC=CC=C5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	412.4 (M + H)	2.85

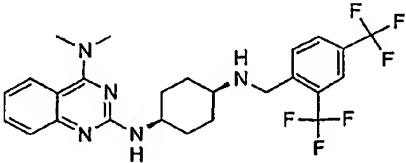
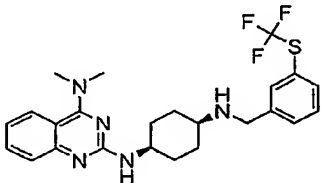
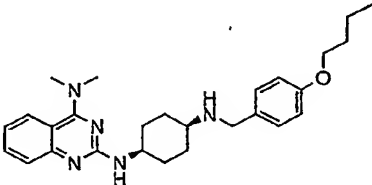
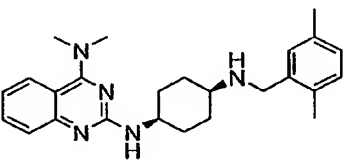
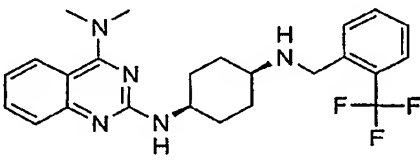
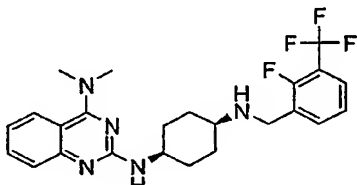
Example No.	Structure	ESI-MS	Retention Time (min)
3305	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4ccc(C(=O)O)cc4)cc5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.67
3306	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4cc(OC)c(N(C)C)cc4)cc5ccccc15</chem> $3\text{CF}_3\text{CO}_2\text{H}$	449.4 (M + H)	2.74
3307	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4ccc(F)cc4)cc5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	394.4 (M + H)	2.86
3308	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4cc(Cl)c(Cl)c(Cl)c4)cc5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.2 (M + H)	3.38
3309	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4cc(F)(F)ccccc4)cc5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	3.09
3310	 <chem>CN(C)c1nc2c(ncn2C1NCC3CCCCC3NCc4ccccc4)cc5ccccc15</chem> $2\text{CF}_3\text{CO}_2\text{H}$	376.4 (M + H)	2.82

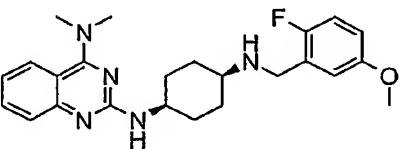
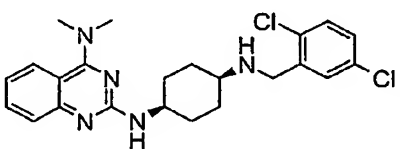
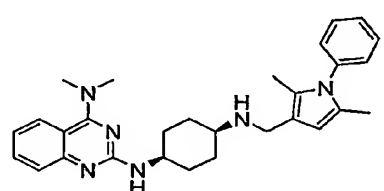
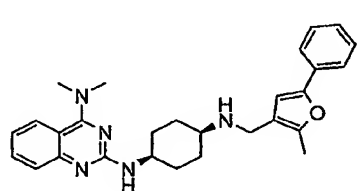
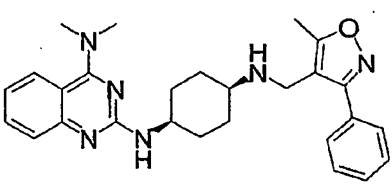
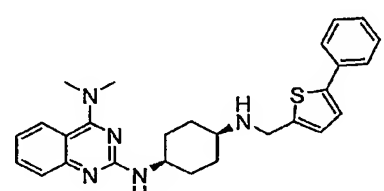
Example No.	Structure	ESI-MS	Retention Time (min)
3311	 2CF ₃ CO ₂ H	406.4 (M + H)	2.87
3312	 2CF ₃ CO ₂ H	436.4 (M + H)	2.91
3313	 2CF ₃ CO ₂ H	426.2 (M + H)	3.13
3314	 2CF ₃ CO ₂ H	436.4 (M + H)	2.99
3315	 2CF ₃ CO ₂ H	454.0 (M + H)	2.97
3316	 2CF ₃ CO ₂ H	412.4 (M + H)	2.92

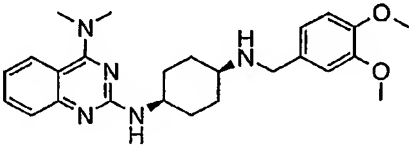
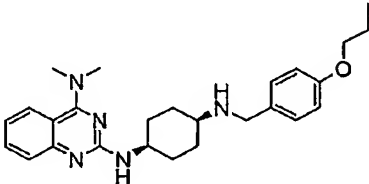
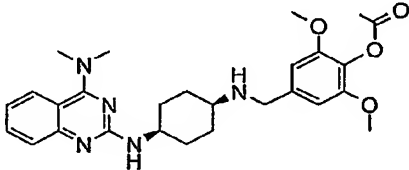
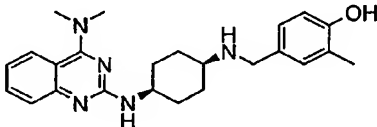
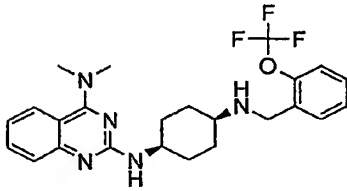
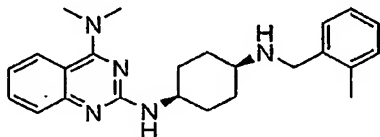
Example No.	Structure	ESI-MS	Retention Time (min)
3317	 $2\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	2.95
3318	 $2\text{CF}_3\text{CO}_2\text{H}$	390.4 (M + H)	2.95
3319	 $2\text{CF}_3\text{CO}_2\text{H}$	396.2 (M + H)	2.89
3320	 $2\text{CF}_3\text{CO}_2\text{H}$	438.2 (M + H)	2.76
3321	 $3\text{CF}_3\text{CO}_2\text{H}$	445.4 (M + H)	3.16
3322	 $3\text{CF}_3\text{CO}_2\text{H}$	415.4 (M + H)	2.96

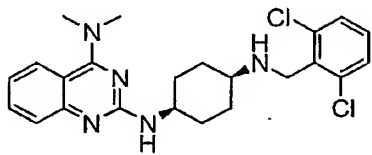
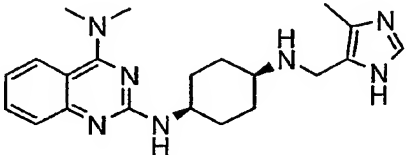
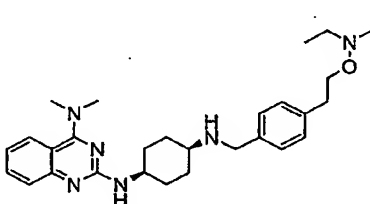
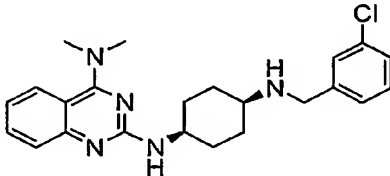
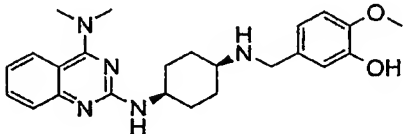
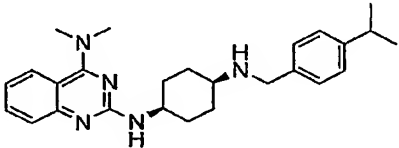
Example No.	Structure	ESI-MS	Retention Time (min)
3323	 $3\text{CF}_3\text{CO}_2\text{H}$	445.4 (M + H)	2.96
3324	 $2\text{CF}_3\text{CO}_2\text{H}$	504.2 (M + H)	3.11
3325	 $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.17
3326	 $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.27
3327	 $2\text{CF}_3\text{CO}_2\text{H}$	514.4 (M + H)	3.07
3328	 $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	2.99

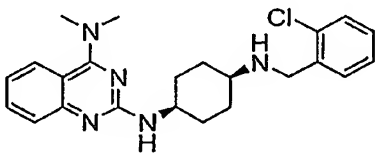
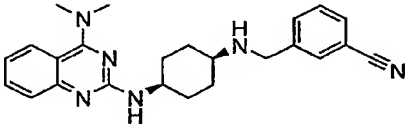
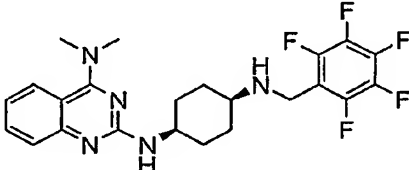
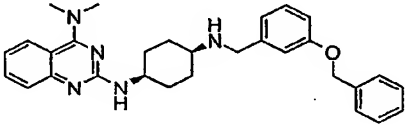
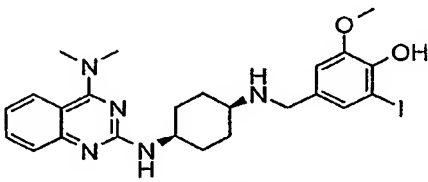
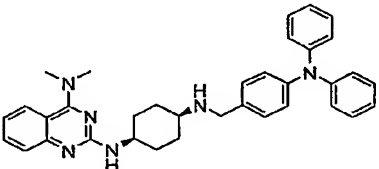
Example No.	Structure	ESI-MS	Retention Time (min)
3329	 $2\text{CF}_3\text{CO}_2\text{H}$	433.2 (M + H)	2.63
3330	 $2\text{CF}_3\text{CO}_2\text{H}$	518.4 (M + H)	3.63
3331	 $2\text{CF}_3\text{CO}_2\text{H}$	500.4 (M + H)	3.09
3332	 $3\text{CF}_3\text{CO}_2\text{H}$	379.4 (M + H)	2.77
3333	 $2\text{CF}_3\text{CO}_2\text{H}$	460.2 (M + H)	3.31
3334	 $2\text{CF}_3\text{CO}_2\text{H}$	512.4 (M + H)	3.51

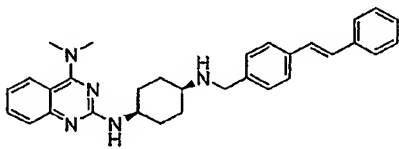
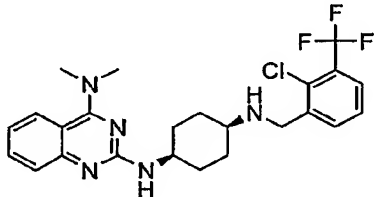
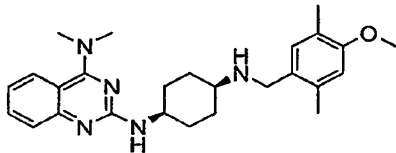
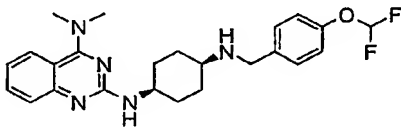
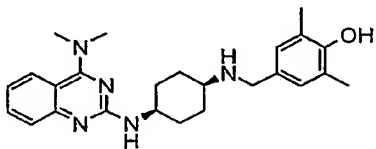
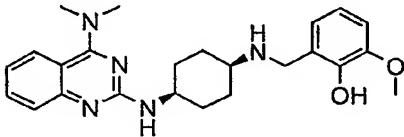
Example No.	Structure	ESI-MS	Retention Time (min)
3335	 2CF ₃ CO ₂ H	512.6 (M + H)	3.51
3336	 2CF ₃ CO ₂ H	476.2 (M + H)	3.39
3337	 2CF ₃ CO ₂ H	448.4 (M + H)	3.42
3338	 2CF ₃ CO ₂ H	404.4 (M + H)	3.17
3339	 2CF ₃ CO ₂ H	444.4 (M + H)	3.13
3340	 2CF ₃ CO ₂ H	462.2 (M + H)	3.21

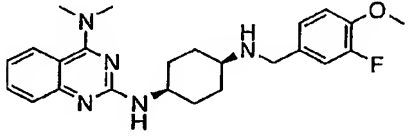
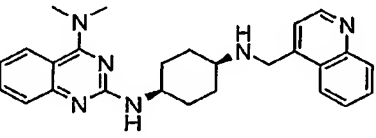
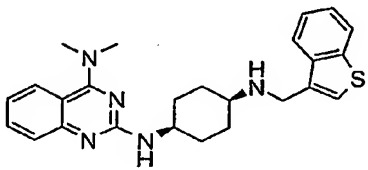
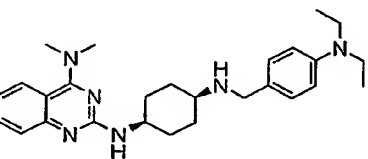
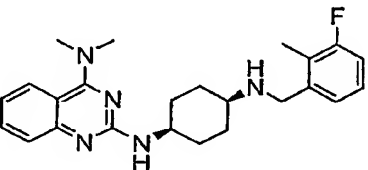
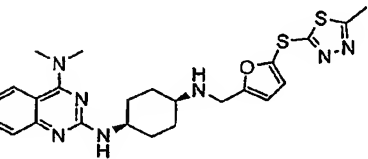
Example No.	Structure	ESI-MS	Retention Time (min)
3341	 $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	2.97
3342	 $2\text{CF}_3\text{CO}_2\text{H}$	444.6 (M + H)	3.16
3343	 $3\text{CF}_3\text{CO}_2\text{H}$	469.4 (M + H)	3.47
3344	 $2\text{CF}_3\text{CO}_2\text{H}$	456.4 (M + H)	3.47
3345	 $2\text{CF}_3\text{CO}_2\text{H}$	457.4 (M + H)	3.09
3346	 $2\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.37

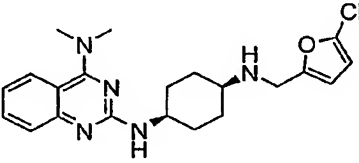
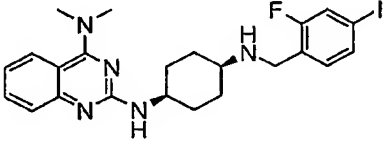
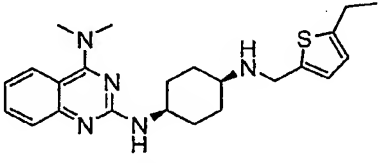
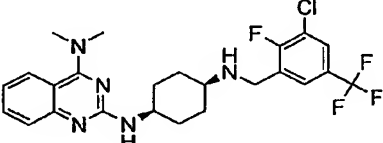
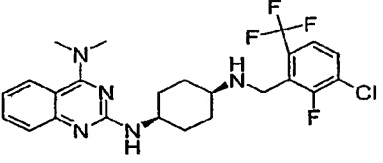
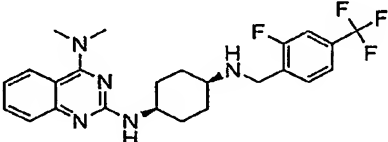
Example No.	Structure	ESI-MS	Retention Time (min)
3347	 <chem>COc1ccc(cc1CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C)CC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	2.83
3348	 <chem>CCOC1=CC=C(C=C1)CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.30
3349	 <chem>CC(=O)OC1=CC(OC)=CC=C1CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	494.4 (M + H)	2.98
3350	 <chem>OC1=CC=C(C=C1)CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	406.4 (M + H)	2.80
3351	 <chem>FC(F)(F)Oc1ccc(cc1)CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	3.20
3352	 <chem>Cc1ccccc1CN[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	390.4 (M + H)	2.97

Example No.	Structure	ESI-MS	Retention Time (min)
3353	 2CF ₃ CO ₂ H	444.2 (M + H)	3.01
3354	 3CF ₃ CO ₂ H	380.2 (M + H)	2.27
3355	 2CF ₃ CO ₂ H	491.4 (M + H)	2.55
3356	 2CF ₃ CO ₂ H	410.4 (M + H)	3.05
3357	 2CF ₃ CO ₂ H	422.2 (M + H)	2.69
3358	 2CF ₃ CO ₂ H	418.6 (M + H)	3.36

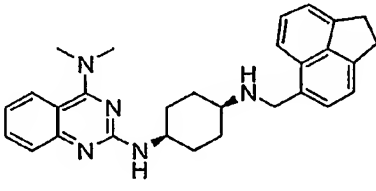
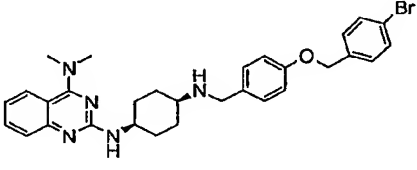
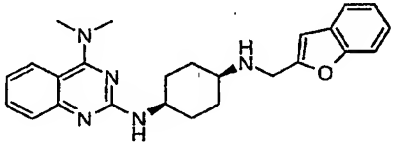
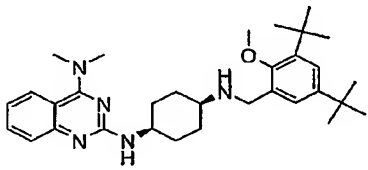
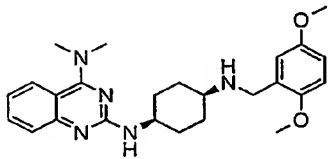
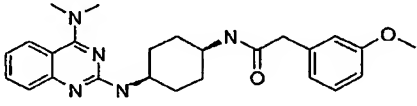
Example No.	Structure	ESI-MS	Retention Time (min)
3359	 $2\text{CF}_3\text{CO}_2\text{H}$	410.4 (M + H)	2.97
3360	 $2\text{CF}_3\text{CO}_2\text{H}$	401.2 (M + H)	2.81
3361	 $2\text{CF}_3\text{CO}_2\text{H}$	466.2 (M + H)	3.01
3362	 $2\text{CF}_3\text{CO}_2\text{H}$	482.4 (M + H)	3.43
3363	 $2\text{CF}_3\text{CO}_2\text{H}$	548.4 (M + H)	3.03
3364	 $3\text{CF}_3\text{CO}_2\text{H}$	543.6 (M + H)	3.95

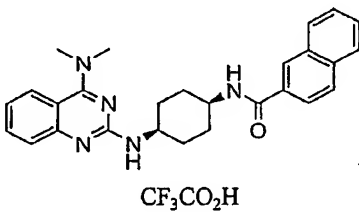
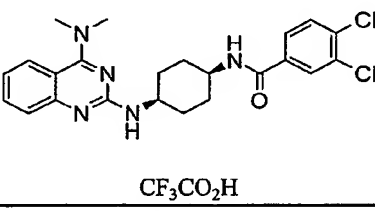
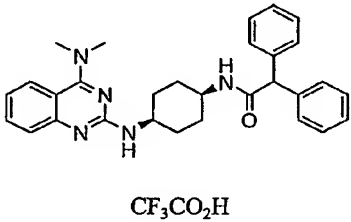
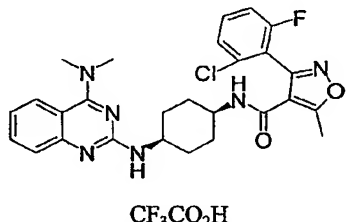
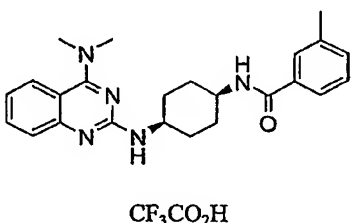
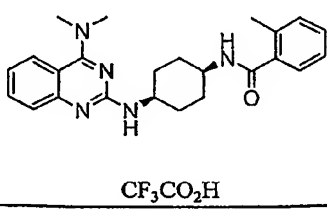
Example No.	Structure	ESI-MS	Retention Time (min)
3365	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	3.64
3366	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	478.4 (M + H)	3.29
3367	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.20
3368	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	442.4 (M + H)	3.09
3369	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.87
3370	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3Nc4ccc(cc4)COC5=CC=CC=C5)nc6ccccc16</chem> $2\text{CF}_3\text{CO}_2\text{H}$	422.2 (M + H)	2.79

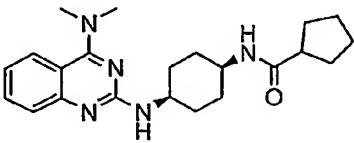
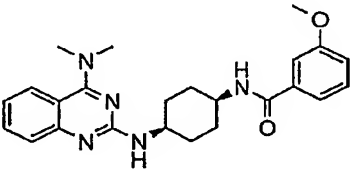
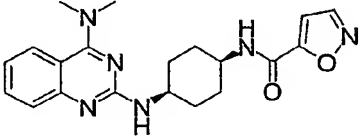
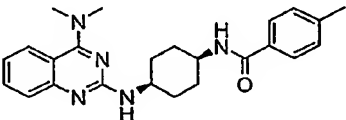
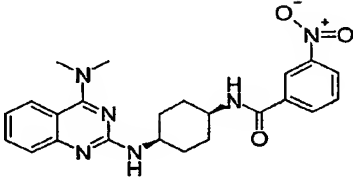
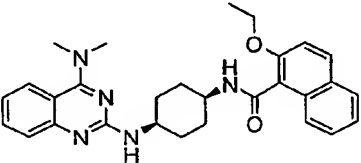
Example No:	Structure	ESI-MS	Retention Time (min)
3371	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3CN(C)C4=CC=C(C=C4)OC(F)=C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	2.96
3372	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3CN(C)C4=CC5=C(NC4=CC=CC5)C6=CC=CC=C6</chem> $3\text{CF}_3\text{CO}_2\text{H}$	427.2 (M + H)	2.53
3373	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3CN(C)C4=CC5=C(C=C4)SC6=CC=CC=C56</chem> $2\text{CF}_3\text{CO}_2\text{H}$	432.4 (M + H)	3.12
3374	 <chem>CCN(CC)Cc1ccc(cc1)CN(C)C2CCCCC2NC3=C(NC4=CC=CC=C4N3C)C5=CC=CC=C5</chem> $3\text{CF}_3\text{CO}_2\text{H}$	447.4 (M + H)	2.45
3375	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3CN(C)C4=CC=C(C=C4)C(F)=C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	408.2 (M + H)	3.02
3376	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3CN(C)C4=CC=CC=C4N3C5C=CC(OC5)SC6=NN=N6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	2.81

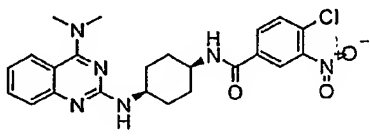
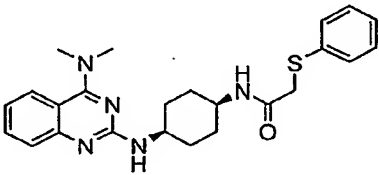
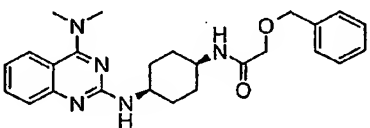
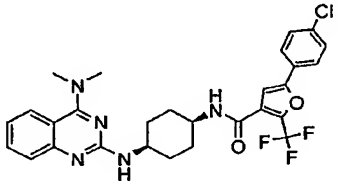
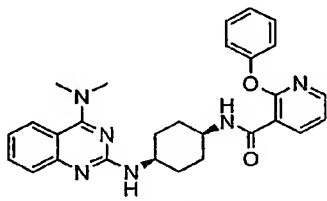
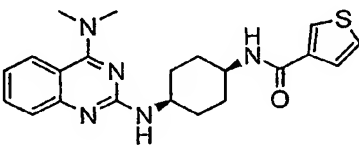
Example No.	Structure	ESI-MS	Retention Time (min)
3377	 $2\text{CF}_3\text{CO}_2\text{H}$	400.2 (M + H)	2.81
3378	 $2\text{CF}_3\text{CO}_2\text{H}$	520.2 (M + H)	3.14
3379	 $2\text{CF}_3\text{CO}_2\text{H}$	410.4 (M + H)	3.12
3380	 $2\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	3.40
3381	 $2\text{CF}_3\text{CO}_2\text{H}$	496.4 (M + H)	3.17
3382	 $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.19

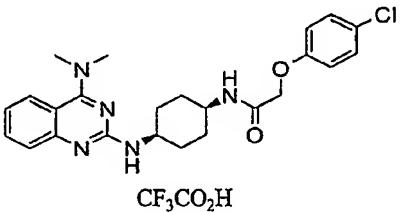
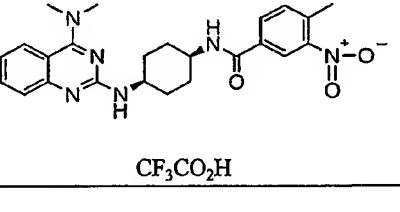
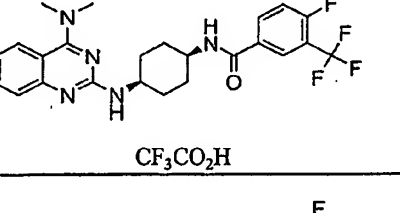
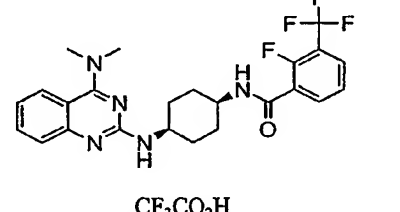
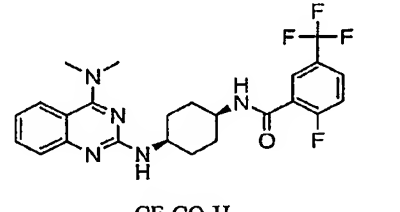
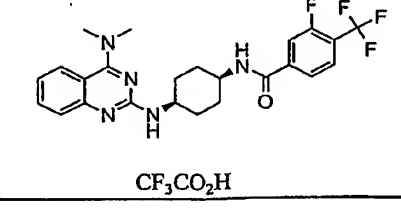
Example No.	Structure	ESI-MS	Retention Time:(min)
3383	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4=CC=C(C=C4)C(F)(F)F)cc1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	3.28
3384	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4=CC(=C(C=C4)OC)C(F)=C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	2.74
3385	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4=CC(=C(C=C4)OC)OC</chem> $2\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	2.89
3386	 <chem>CC(C1=CC=CC=C1)NC2CCCCC2NC3=CC4=CC=CC=C4N(C)N3</chem> $2\text{CF}_3\text{CO}_2\text{H}$	404.4 (M + H)	3.09
3387	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4=CC5=C(C=C4)OC(=C5)C(F)=CC=C5Cl</chem> $2\text{CF}_3\text{CO}_2\text{H}$	482.2 (M + H)	3.29
3388	 <chem>CN(C)c1nc2c(ncn2C3CCCCC3NC4=CC=CC=C4NC5CCN(C)C5</chem> $3\text{CF}_3\text{CO}_2\text{H}$	458.4 (M + H)	2.99

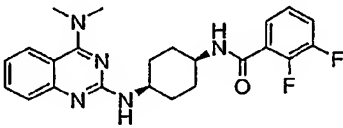
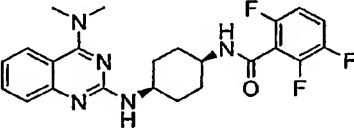
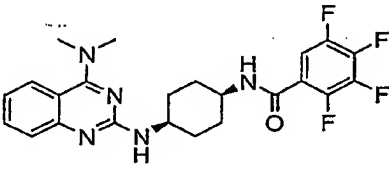
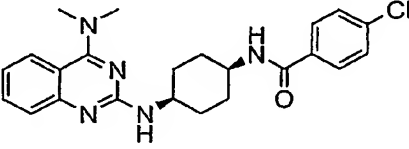
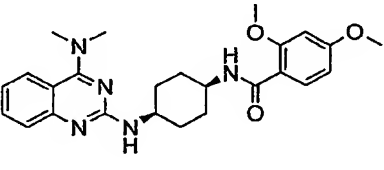
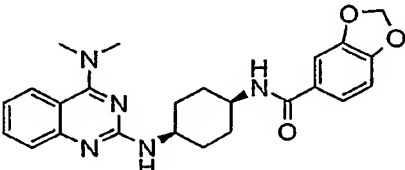
Example No.	Structure	ESI-MS	Retention Time (min)
3389	 $2\text{CF}_3\text{CO}_2\text{H}$	452.2 (M + H)	3.40
3390	 $2\text{CF}_3\text{CO}_2\text{H}$	560.2 (M + H)	3.73
3391	 $2\text{CF}_3\text{CO}_2\text{H}$	416.4 (M + H)	2.99
3392	 $2\text{CF}_3\text{CO}_2\text{H}$	518.6 (M + H)	4.08
3393	 $2\text{CF}_3\text{CO}_2\text{H}$	436.4 (M + H)	2.95
3394	 $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.30

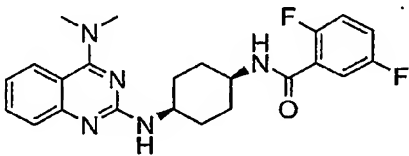
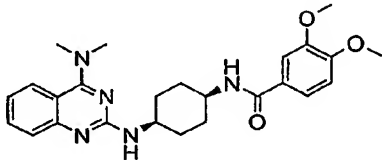
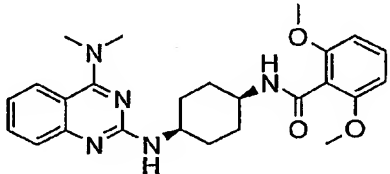
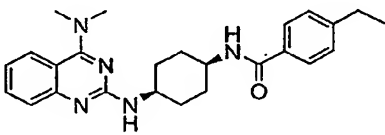
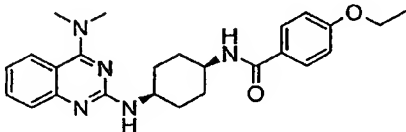
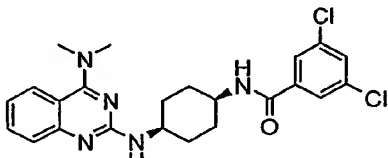
Example No.	Structure	ESI-MS	Retention Time (min)
3395	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)c5ccc6ccccc6c5.CF3C(=O)O</chem>	440.4 (M + H)	4.26
3396	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)c5cc(Cl)cc(Cl)c5.CF3C(=O)O</chem>	458.2 (M + H)	4.39
3397	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)C(c5ccccc5)c6ccccc6.CF3C(=O)O</chem>	480.4 (M + H)	4.37
3398	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)c5cc(F)cc(Cl)c5.CF3C(=O)O</chem>	523.6 (M + H)	4.15
3399	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)c5cccc(C)c5.CF3C(=O)O</chem>	404.4 (M + H)	3.46
3400	 <chem>CC1=C(C)N2C(=N1)c3ccccc3N2C4CCCCC4NC(=O)c5ccccc5C.CF3C(=O)O</chem>	404.4 (M + H)	3.75

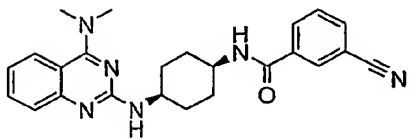
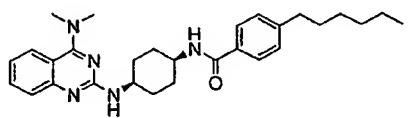
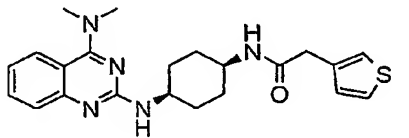
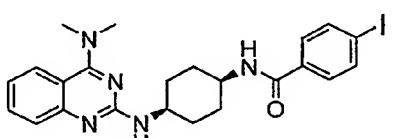
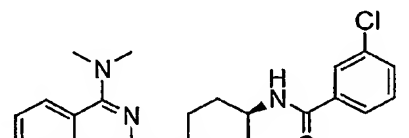
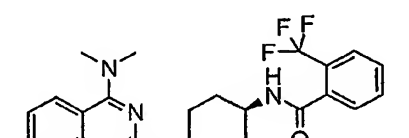
Example No.	Structure	ESI-MS	Retention Time (min)
3401	 CF ₃ CO ₂ H	382.4 (M + H)	3.65
3402	 CF ₃ CO ₂ H	420.4 (M + H)	3.81
3403	 CF ₃ CO ₂ H	381.2 (M + H)	3.33
3404	 CF ₃ CO ₂ H	404.4 (M + H)	3.93
3405	 CF ₃ CO ₂ H	435.2 (M + H)	3.40
3406	 CF ₃ CO ₂ H	484.4 (M + H)	4.15

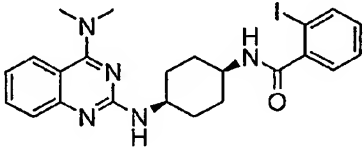
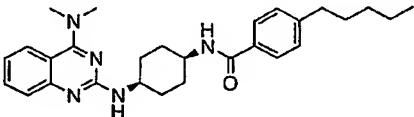
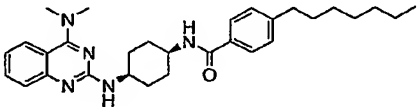
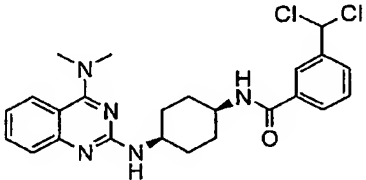
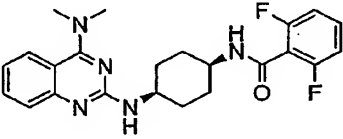
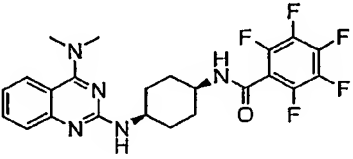
Example No.	Structure	ESI-MS	Retention Time (min)
3407	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)C(=C(C=C3)Cl)[N+](=O)[O-]</chem> $\text{CF}_3\text{CO}_2\text{H}$	469.4 (M + H)	4.20
3408	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)CS4=CC=CC=C4</chem> $\text{CF}_3\text{CO}_2\text{H}$	436.2 (M + H)	3.88
3409	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)OCC4=CC=CC=C4</chem> $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	3.91
3410	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)C4=CC(=C(C=C4)OC(F)(F)F)C5=CC=C(C=C5)Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	558.4 (M + H)	4.92
3411	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)C4=CC=CC=C4OC5=CC=CC=C5N6=CC=CC=C6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	483.4 (M + H)	4.08
3412	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3N3C(=O)C4=CC=C(C=C4)S</chem> $\text{CF}_3\text{CO}_2\text{H}$	396.2 (M + H)	3.68

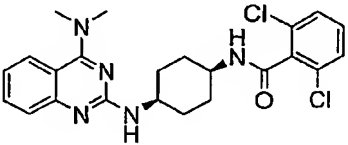
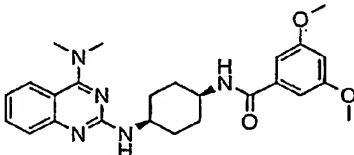
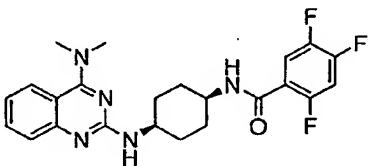
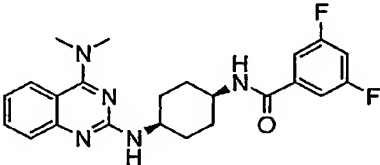
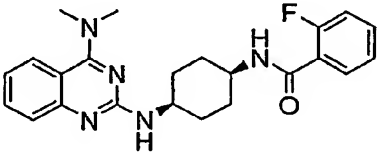
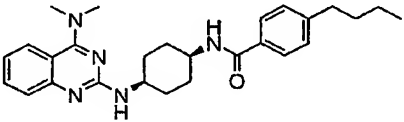
Example No.	Structure	ESI-MS	Retention Time (min)
3413	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)Oc5ccc(Cl)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	454.2 (M + H)	3.70
3414	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)c5ccc([N+](=O)[O-])cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	449.4 (M + H)	4.09
3415	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)c5ccc(C(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	4.33
3416	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)c5cc(F)c(C(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	3.60
3417	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)c5cc(F)c(C(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	4.23
3418	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3C(=O)N[C@H]4CCCC[C@H]4C(=O)c5cc(F)c(C(F)(F)F)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.4 (M + H)	4.38

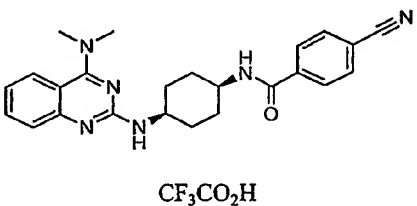
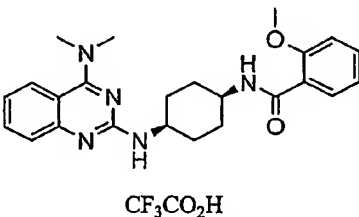
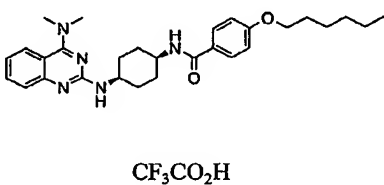
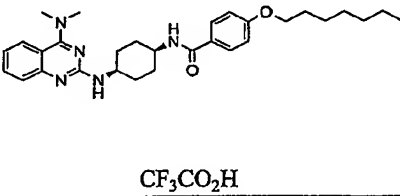
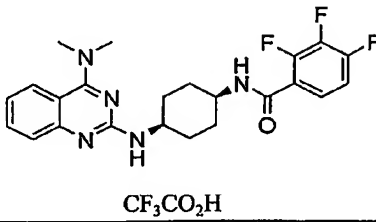
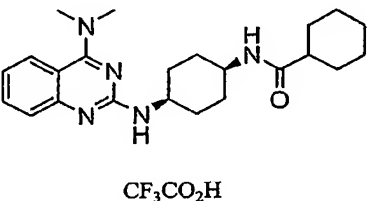
Example No.	Structure	ESI-MS	Retention Time (min)
3419	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4cc(F)c(F)cc4</chem> <chem>CF3CO2H</chem>	426.2 (M + H)	3.87
3420	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4cc(F)c(F)c(F)c4</chem> <chem>CF3CO2H</chem>	444.4 (M + H)	3.86
3421	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4cc(F)c(F)c(F)c4F</chem> <chem>CF3CO2H</chem>	462.2 (M + H)	4.15
3422	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4ccc(Cl)cc4</chem> <chem>CF3CO2H</chem>	424.2 (M + H)	4.06
3423	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4cc(OC)c(OC)cc4</chem> <chem>CF3CO2H</chem>	450.4 (M + H)	4.03
3424	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N3CCCC[C@H]3NC(=O)c4cc5c(c(c4)O)OCC5</chem> <chem>CF3CO2H</chem>	434.2 (M + H)	3.75

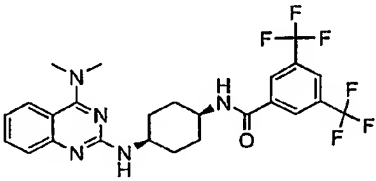
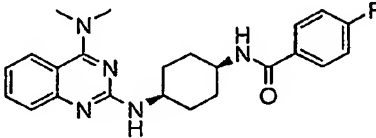
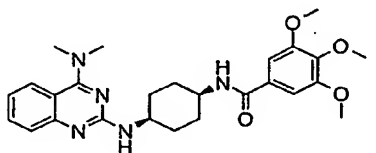
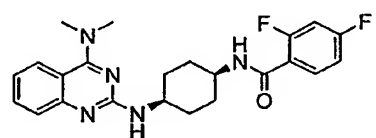
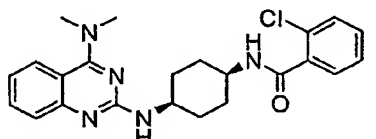
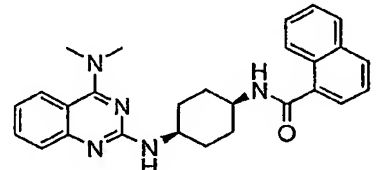
Example No.	Structure	ESI-MS	Retention Time (min)
3425	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4cc(F)cc(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	3.88
3426	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4cc(OC)c(OC)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	3.64
3427	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4cc(OC)c(OC)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	3.55
3428	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4ccc(CC)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	418.6 (M + H)	4.17
3429	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4ccc(OCC)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	4.03
3430	 <chem>CC1=NC2=CC=CC=C2N1C(=N)N[C@H]3CCCC[C@H]3NC(=O)c4cc(Cl)cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	4.45

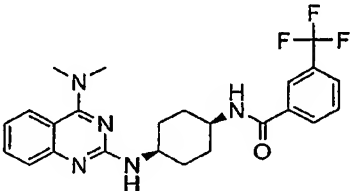
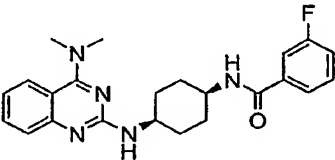
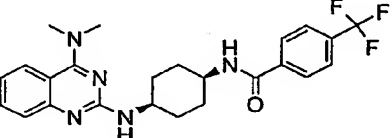
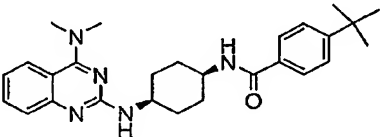
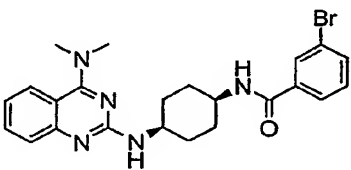
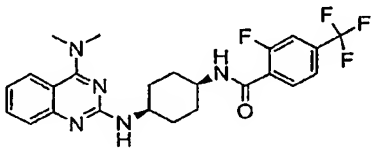
Example No.	Structure	ESI-MS	Retention Time (min)
3431	 <chem>CC1=NC2=CC=CC=C2N1N=C(NC3CCCCC3C(=O)NC4=CC=CC=C4C#N)C5=CC=CC=C5</chem> $\text{CF}_3\text{CO}_2\text{H}$	415.4 (M + H)	3.76
3432	 <chem>CCCCCc1ccc(cc1)C(=O)NC2CCCCC2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	5.06
3433	 <chem>Cc1ccsc1C(=O)NC2CCCCC2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	410.2 (M + H)	3.64
3434	 <chem>Ic1ccc(cc1)C(=O)NC2CCCCC2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.2 (M + H)	4.24
3435	 <chem>Clc1ccc(cc1)C(=O)NC2CCCCC2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	4.09
3436	 <chem>Cc1ccccc1C(=O)NC2CCCCC2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.89

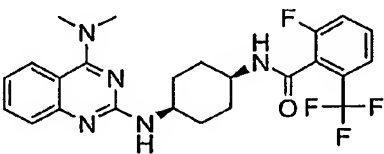
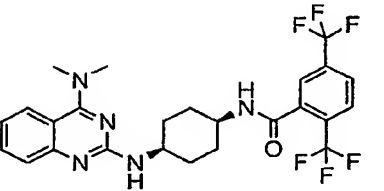
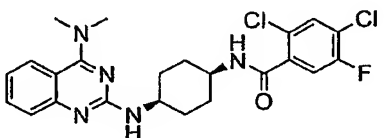
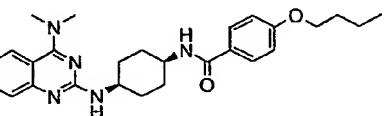
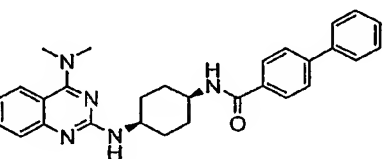
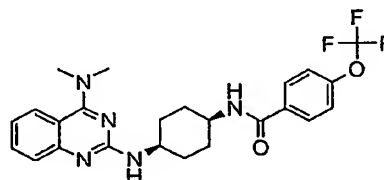
Example No.	Structure	ESI-MS	Retention Time (min)
3437	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	516.2 (M + H)	3.88
3438	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	4.86
3439	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	488.4 (M + H)	4.70
3440	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	472.4 (M + H)	4.29
3441	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	3.69
3442	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC=CC=C3NC4=CC=CC=C4C(=O)N[C@H]5CCCC[C@H]5N</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.2 (M + H)	4.16

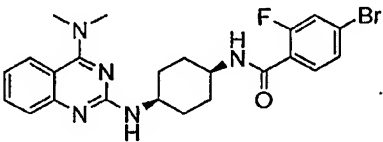
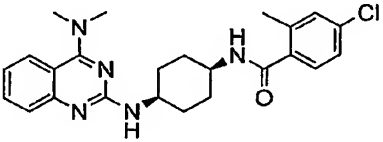
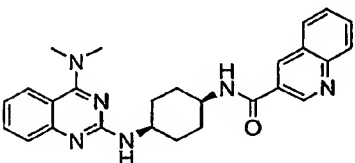
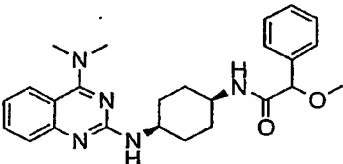
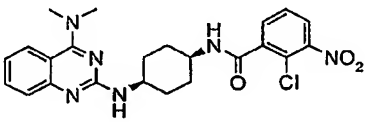
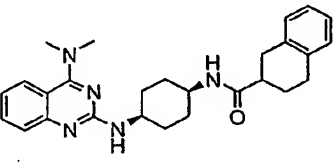
Example No.	Structure	ESI-MS	Retention Time (min)
3443	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5cc(Cl)cc(Cl)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	458.2 (M + H)	3.91
3444	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5cc(OC)c(OC)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	3.95
3445	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5cc(F)c(F)c(F)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	4.01
3446	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5cc(F)c(F)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	4.00
3447	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5ccccc5F</chem> $\text{CF}_3\text{CO}_2\text{H}$	408.4 (M + H)	3.75
3448	 <chem>CC1=NC2=CC=CC=C2N1C3=CC=CC=C3N3C4CCCCC4NC(=O)c5ccc(CCC)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	446.6 (M + H)	4.65

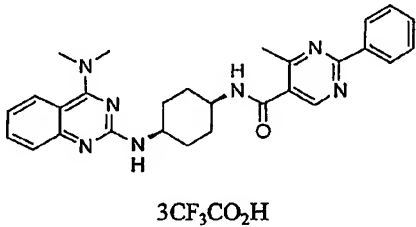
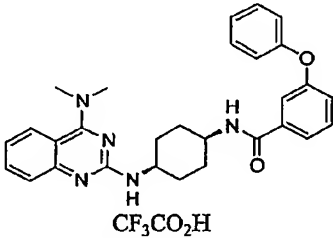
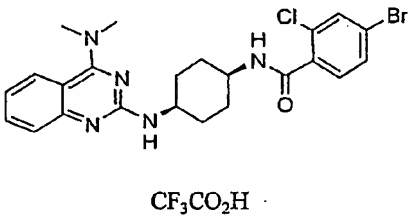
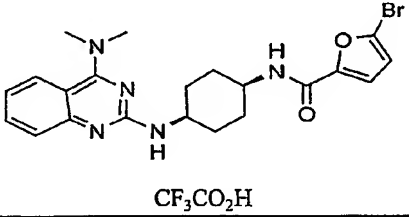
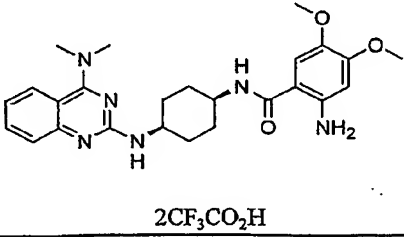
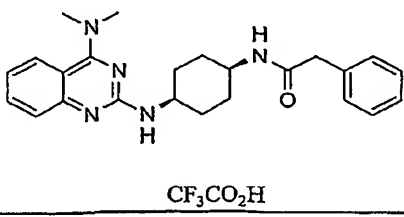
Example No.	Structure	ESI-MS	Retention Time (min)
3449	 CF ₃ CO ₂ H	415.2 (M + H)	3.75
3450	 CF ₃ CO ₂ H	420.4 (M + H)	3.91
3451	 CF ₃ CO ₂ H	490.4 (M + H)	4.99
3452	 CF ₃ CO ₂ H	504.4 (M + H)	5.16
3453	 CF ₃ CO ₂ H	444.4 (M + H)	4.00
3454	 CF ₃ CO ₂ H	396.2 (M + H)	3.85

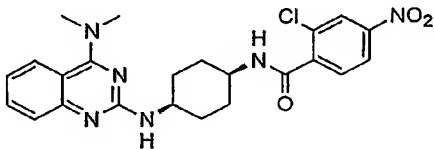
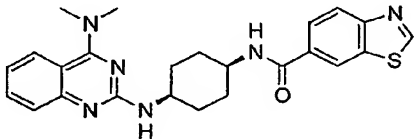
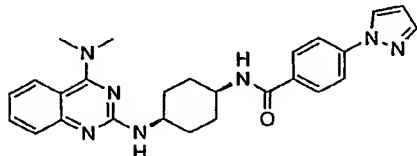
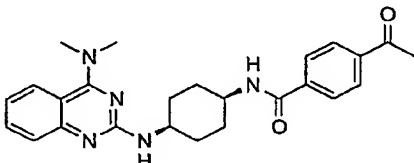
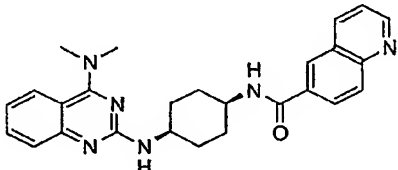
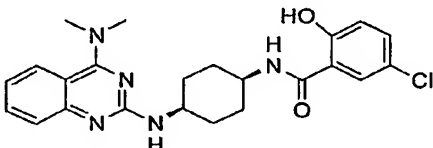
Example No.	Structure	ESI-MS	Retention Time (min)
3455	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=C(C4)C(F)(F)F)C(F)(F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	526.6 (M + H)	4.69
3456	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=CC=C4)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	408.4 (M + H)	3.30
3457	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=C(C4)OC)OC</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.4 (M + H)	3.76
3458	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=C(C4)F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	426.2 (M + H)	3.86
3459	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=CC=C4)Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	424.2 (M + H)	3.76
3460	 <chem>CC1=NC2=CC=CC=C2N(C)N1C3=CC(=CC=C3)NC(=O)N4C=CC(=CC=C4C5=CC=CC=C5)</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.4 (M + H)	4.05

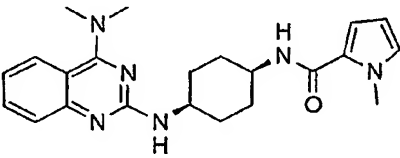
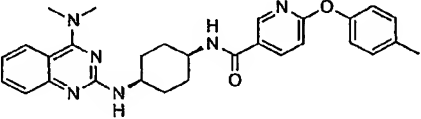
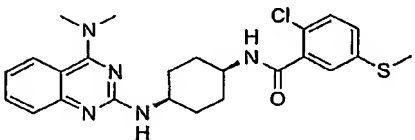
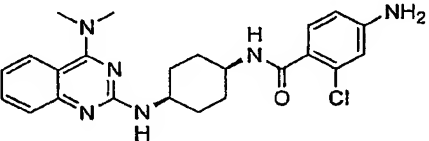
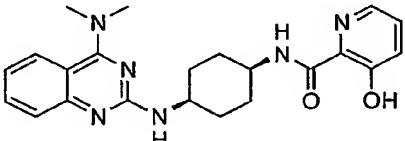
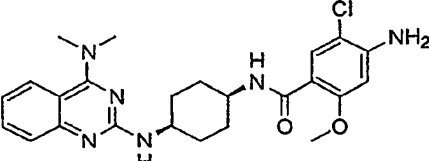
Example No.	Structure	ESI-MS	Retention Time (min)
3461	 CF ₃ CO ₂ H	458.4 (M + H)	4.25
3462	 CF ₃ CO ₂ H	408.2 (M + H)	3.84
3463	 CF ₃ CO ₂ H	458.2 (M + H)	4.25
3464	 CF ₃ CO ₂ H	446.6 (M + H)	4.44
3465	 CF ₃ CO ₂ H	470.2 (M + H)	4.13
3466	 CF ₃ CO ₂ H	476.2 (M + H)	4.25

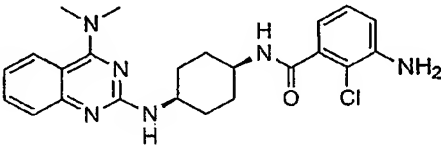
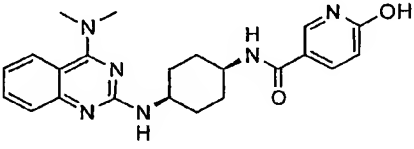
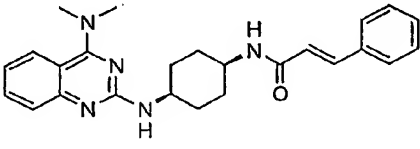
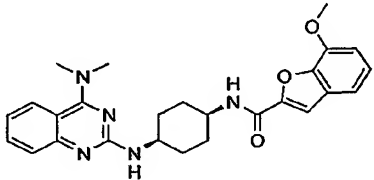
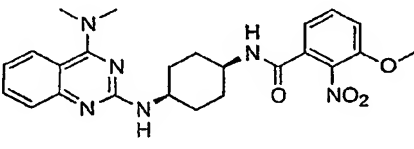
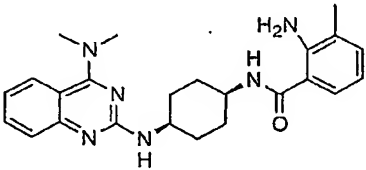
Example No.	Structure	ESI-MS	Retention Time (min)
3467	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1cc(F)cc(C(F)(F)F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.92
3468	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1cc(F)c(C(F)F)c(F)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	526.4 (M + H)	4.31
3469	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1cc(F)cc(Cl)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	4.15
3470	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1ccc(OCC)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	462.2 (M + H)	4.48
3471	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1ccc(cc1)-c2ccccc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	4.45
3472	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c1ccc(OC(F)(F)F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	4.29

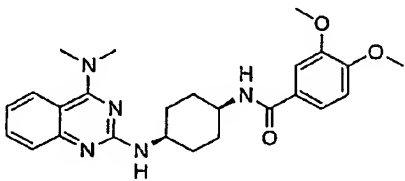
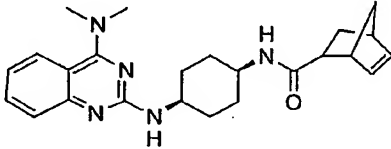
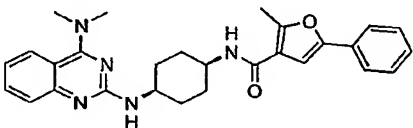
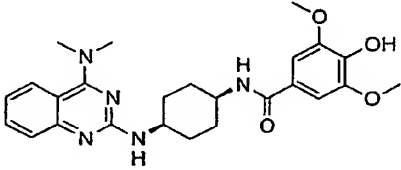
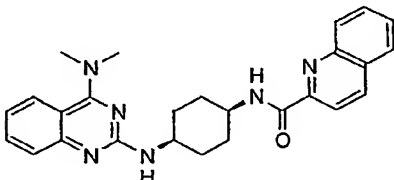
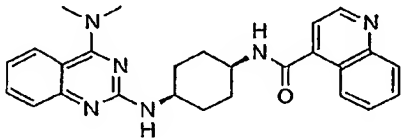
Example No.	Structure	ESI-MS	Retention Time (min)
3473	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c4cc(Br)cc(F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	486.2 (M + H)	4.32
3474	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c4cc(Cl)cc(C)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	438.4 (M + H)	4.31
3475	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c4cnc5ccccc5c4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	441.4 (M + H)	3.75
3476	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)C(c4ccccc4)OC</chem> $\text{CF}_3\text{CO}_2\text{H}$	434.4 (M + H)	4.10
3477	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c4cc([N+](=O)[O-])cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	469.4 (M + H)	4.19
3478	 <chem>CC1=NC2=CC=CC=C2N(C)N1[C@H]3CCCC[C@H]3NC(=O)c4c5ccccc5cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	444.4 (M + H)	4.36

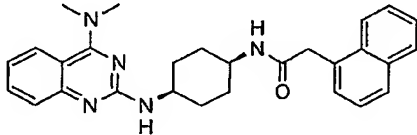
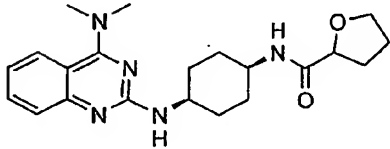
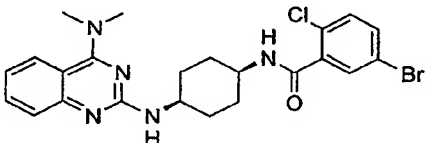
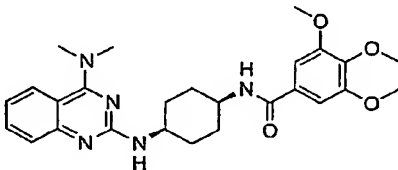
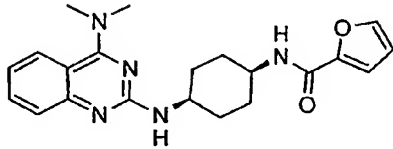
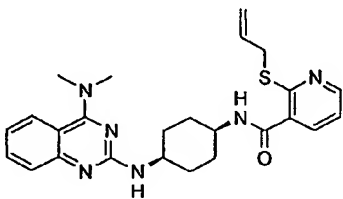
Example No.	Structure	ESI-MS	Retention Time (min)
3479	 3CF ₃ CO ₂ H	482.4 (M + H)	4.35
3480	 CF ₃ CO ₂ H	482.4 (M + H)	4.64
3481	 CF ₃ CO ₂ H	502.2 (M + H)	4.37
3482	 CF ₃ CO ₂ H	458.2 (M + H)	4.08
3483	 2CF ₃ CO ₂ H	465.4 (M + H)	3.66
3484	 CF ₃ CO ₂ H	404.4 (M + H)	4.03

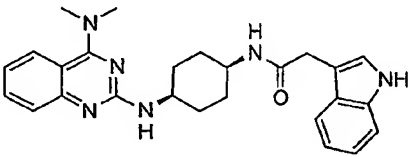
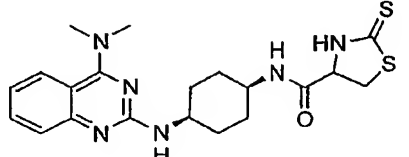
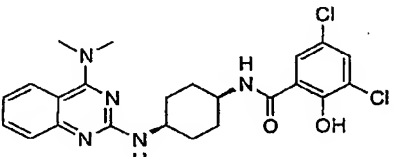
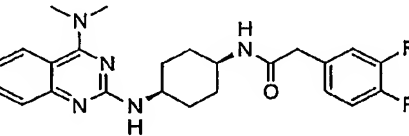
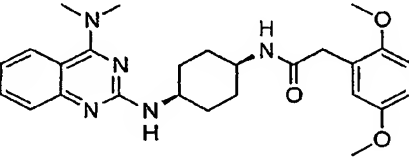
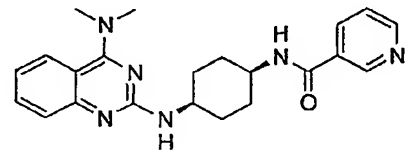
Example No.	Structure	ESI-MS	Retention Time (min)
3485	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4cc([N+](=O)[O-])cc(Cl)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	469.4 (M + H)	4.23
3486	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4cc5ncsc5cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	447.4 (M + H)	3.94
3487	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4ccc5c(c4)nn5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	456.2 (M + H)	4.07
3488	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4ccc(C(C)=O)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	432.4 (M + H)	3.99
3489	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4ccc5c(c4)nc6ccccc65</chem> $2\text{CF}_3\text{CO}_2\text{H}$	441.3 (M + H)	1.70
3490	 <chem>CC1=NC2=C(N1)N=CN=C2N[C@H]3CCCC[C@H]3NC(=O)c4cc(Cl)cc(O)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.2 (M + H)	4.57

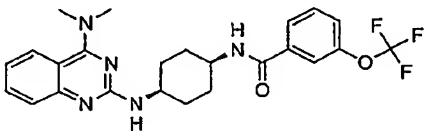
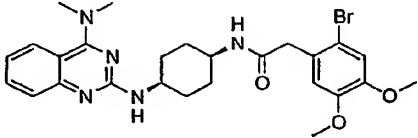
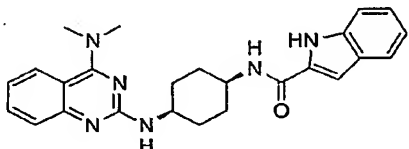
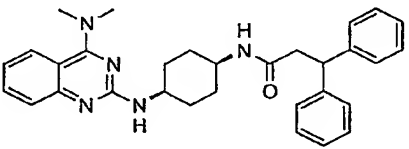
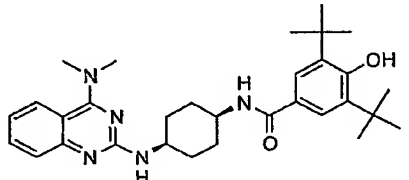
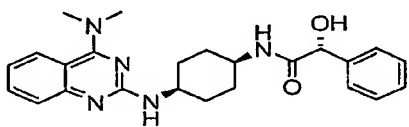
Example No.	Structure	ESI-MS	Retention Time (min)
3491	 $2\text{CF}_3\text{CO}_2\text{H}$	393.4 (M + H)	4.01
3492	 $2\text{CF}_3\text{CO}_2\text{H}$	497.4 (M + H)	4.45
3493	 $\text{CF}_3\text{CO}_2\text{H}$	470.2 (M + H)	2.40
3494	 $2\text{CF}_3\text{CO}_2\text{H}$	439.4 (M + H)	1.92
3495	 $2\text{CF}_3\text{CO}_2\text{H}$	407.4 (M + H)	2.30
3496	 $2\text{CF}_3\text{CO}_2\text{H}$	469.5 (M + H)	2.27

Example No.	Structure	ESI-MS	Retention Time (min)
3497	 $2\text{CF}_3\text{CO}_2\text{H}$	439.4 (M + H)	1.93
3498	 $2\text{CF}_3\text{CO}_2\text{H}$	407.4 (M + H)	1.62
3499	 $\text{CF}_3\text{CO}_2\text{H}$	416.3 (M + H)	2.34
3500	 $\text{CF}_3\text{CO}_2\text{H}$	460.4 (M + H)	2.46
3501	 $\text{CF}_3\text{CO}_2\text{H}$	465.4 (M + H)	4.13
3502	 $2\text{CF}_3\text{CO}_2\text{H}$	419.4 (M + H)	3.87

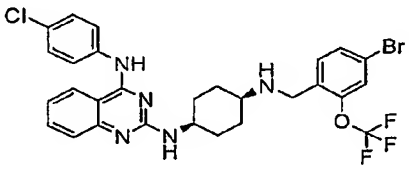
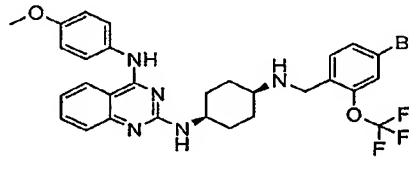
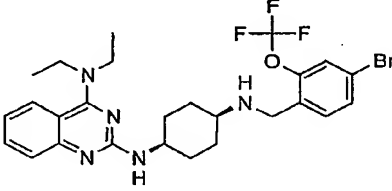
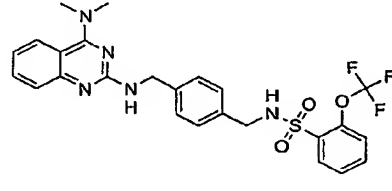
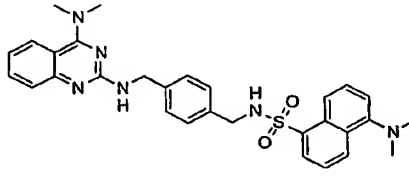
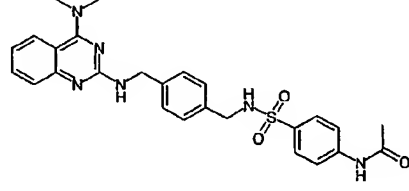
Example No.	Structure	ESI-MS	Retention Time (min)
3503	 <chem>COc1ccc(cc1)C(=O)Nc2ccc(cc2)Nc3ccc(cc3)N(C)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	3.97
3504	 <chem>C1CC2C(C1)C(=O)Nc3ccc(cc3)Nc4ccc(cc4)N(C)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	406.2 (M + H)	2.18
3505	 <chem>c1ccc(cc1)-c2cc3c(c2)oc(=O)Nc4ccc(cc4)Nc5ccc(cc5)N(C)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	470.4 (M + H)	4.74
3506	 <chem>COc1cc(O)cc(OC)c1C(=O)Nc2ccc(cc2)Nc3ccc(cc3)N(C)C</chem> $\text{CF}_3\text{CO}_2\text{H}$	466.4 (M + H)	3.83
3507	 <chem>c1ccc2c(c1)nc3ccc(cc3n2)C(=O)Nc4ccc(cc4)Nc5ccc(cc5)N(C)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	441.2 (M + H)	4.38
3508	 <chem>c1ccc2c(c1)nc3ccc(cc3n2)C(=O)Nc4ccc(cc4)Nc5ccc(cc5)N(C)C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	441.2 (M + H)	3.62

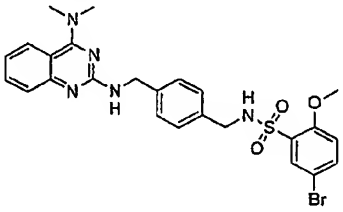
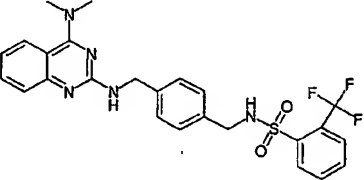
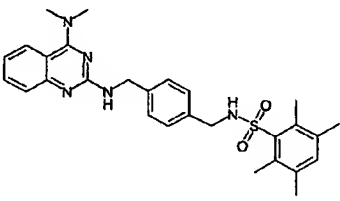
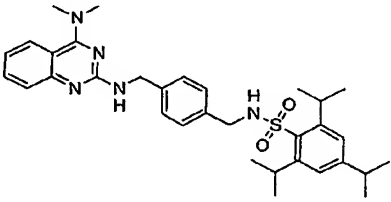
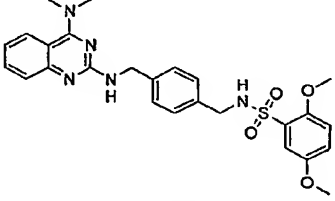
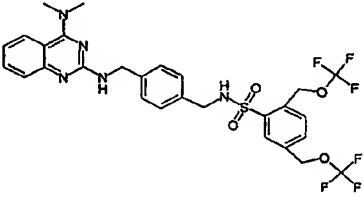
Example No.	Structure	ESI-MS	Retention Time (min)
3509	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5=CC=CC=C5</chem> $\text{CF}_3\text{CO}_2\text{H}$	454.5 (M + H)	2.44
3510	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5OCCC5</chem> $\text{CF}_3\text{CO}_2\text{H}$	384.4 (M + H)	3.67
3511	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5=CC(=C(C=C5)Br)C(=C)Cl</chem> $\text{CF}_3\text{CO}_2\text{H}$	502.2 (M + H)	4.37
3512	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5=CC(=C(C=C5)OC)OC</chem> $\text{CF}_3\text{CO}_2\text{H}$	480.5 (M + H)	2.18
3513	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5=CC=CC=C5O</chem> $\text{CF}_3\text{CO}_2\text{H}$	380.2 (M + H)	3.81
3514	 <chem>CC1(C)N2C=NC3=CC=CC=C3N=C2N1C4CCCCC4NC(=O)CC5=CC=CC=C5S6C=CC=CN6</chem> $2\text{CF}_3\text{CO}_2\text{H}$	463.2 (M + H)	4.23

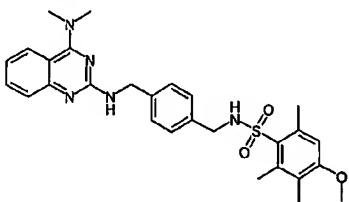
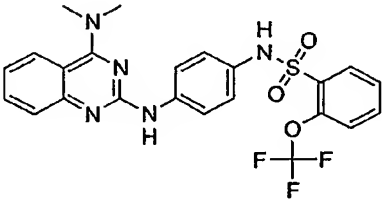
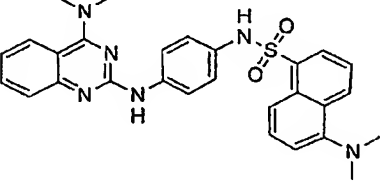
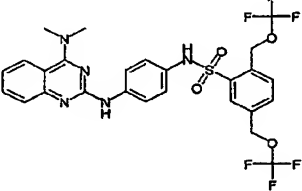
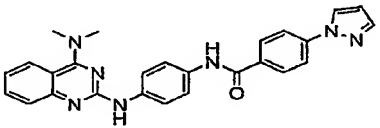
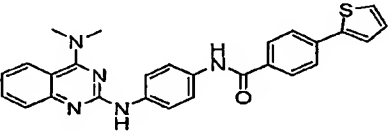
Example No.	Structure	ESI-MS	Retention Time (min)
3515	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)CC4C=CC=C5C(=C4)NCC5</chem> $2\text{CF}_3\text{CO}_2\text{H}$	443.4 (M + H)	2.12
3516	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)CC4=NC(=S)NC=S4</chem> $\text{CF}_3\text{CO}_2\text{H}$	431.1 (M + H)	1.90
3517	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)Cc1cc(Cl)c(Cl)cc1O</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	5.05
3518	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)Cc1cc(F)c(F)cc1</chem> $\text{CF}_3\text{CO}_2\text{H}$	440.5 (M + H)	2.33
3519	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)Cc1cc(OC)cc(OC)c1</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.5 (M + H)	2.20
3520	 <chem>CN1C=NC2=C(N1)C=CC=C2C3CCCCC3NC(=O)Cc1ccccn1</chem> $2\text{CF}_3\text{CO}_2\text{H}$	391.1 (M + H)	1.59

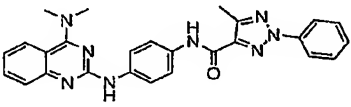
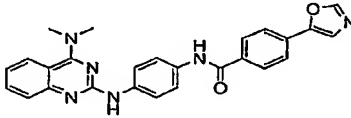
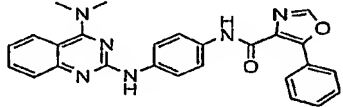
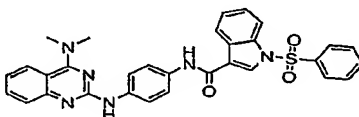
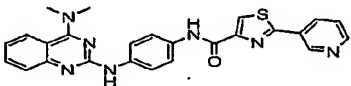
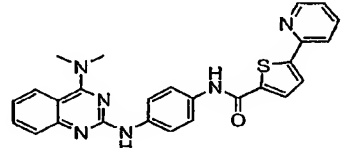
Example No.	Structure	ESI-MS	Retention Time (min)
3521	 <chem>CC1=NC2=CC=CC=C2N(C)C1N[C@H]3CCCC[C@H]3NC(=O)c4ccc(OC(F)(F)F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	474.4 (M + H)	4.53
3522	 <chem>COc1cc(OC)c(Br)cc1NC(=O)[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	542.2 (M + H)	2.26
3523	 <chem>Cc1ccc2c(c1)c(c[nH]2)C(=O)N[C@H]3CCCC[C@H]3Nc4nc5ccccc5n4C</chem> $2\text{CF}_3\text{CO}_2\text{H}$	429.3 (M + H)	2.41
3524	 <chem>c1ccccc1C(Cc2ccccc2)C(=O)N[C@H]3CCCC[C@H]3Nc4nc5ccccc5n4C</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.6 (M + H)	2.59
3525	 <chem>CC(C)(C)c1cc(O)cc(C(C)(C)C)c1C(=O)N[C@H]2CCCC[C@H]2Nc3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	518.5 (M + H)	2.96
3526	 <chem>CC1=NC2=CC=CC=C2N(C)C1N[C@H]3CCCC[C@H]3NC(=O)[C@H](O)c4ccccc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	420.4 (M + H)	2.19

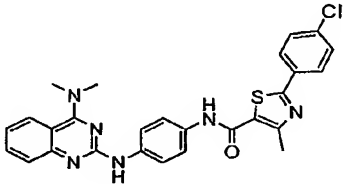
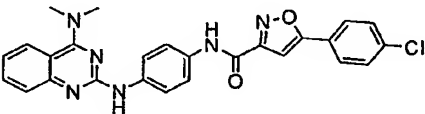
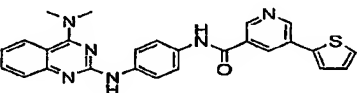
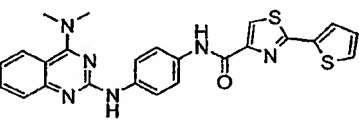
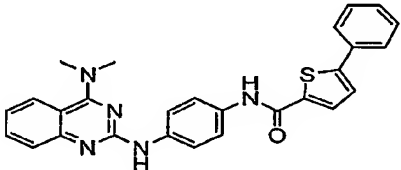
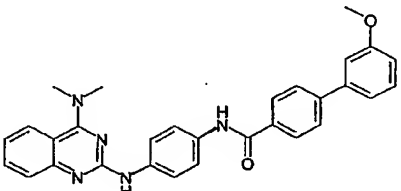
Example No.	Structure	ESI-MS	Retention Time (min)
3527	 CF ₃ CO ₂ H	420.4 (M + H)	2.19
3528	 2CF ₃ CO ₂ H	552.0 (M + H)	2.45
3529	 2CF ₃ CO ₂ H	564.2 (M + H)	2.48
3530	 2CF ₃ CO ₂ H	606.0 (M + H)	2.86
3531	 2CF ₃ CO ₂ H	586.2 (M + H)	3.20
3532	 2CF ₃ CO ₂ H	614.4 (M + H)	2.76

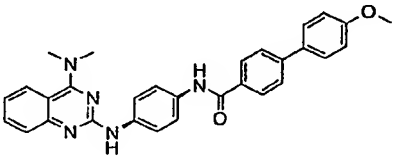
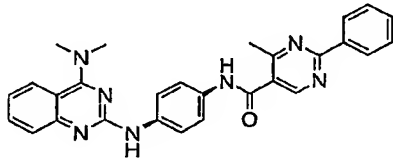
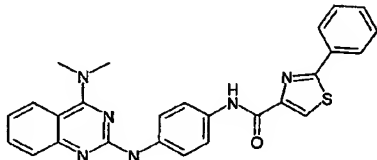
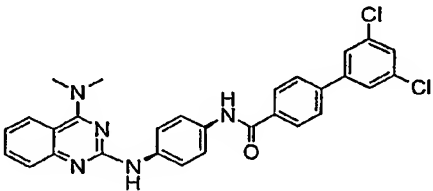
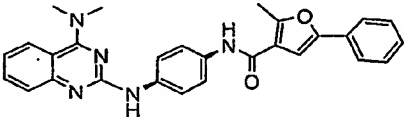
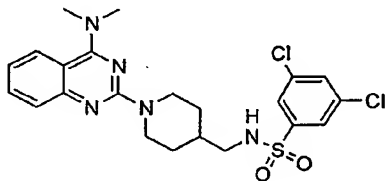
Example No.	Structure	ESI-MS	Retention Time (min)
3533	 <p>2CF₃CO₂H</p>	620.0 (M + H)	2.68
3534	 <p>2CF₃CO₂H</p>	616.0 (M + H)	2.56
3535	 <p>2CF₃CO₂H</p>	566.0 (M + H)	2.54
3536	 <p>CF₃CO₂H</p>	532.2 (M + H)	3.35
3537	 <p>2CF₃CO₂H</p>	541.4 (M + H)	3.11
3538	 <p>CF₃CO₂H</p>	505.2 (M + H)	2.98

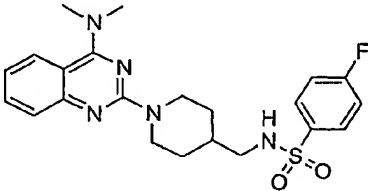
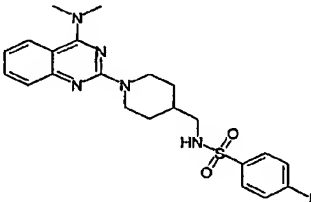
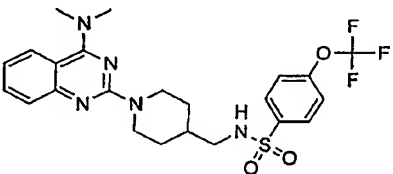
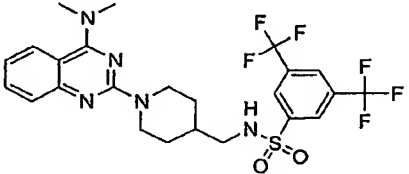
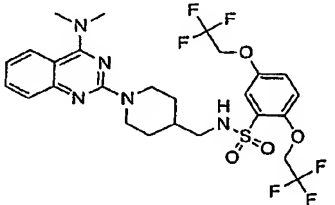
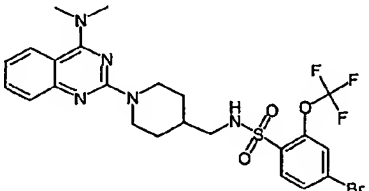
Example No.	Structure	ESI-MS	Retention Time (min)
3539	 $\text{CF}_3\text{CO}_2\text{H}$	556 (M + H)	3.37
3540	 $\text{CF}_3\text{CO}_2\text{H}$	516.4 (M + H)	3.39
3541	 $\text{CF}_3\text{CO}_2\text{H}$	504.4 (M + H)	3.61
3542	 $\text{CF}_3\text{CO}_2\text{H}$	574.4 (M + H)	4.27
3543	 $\text{CF}_3\text{CO}_2\text{H}$	508.2 (M + H)	3.17
3544	 $\text{CF}_3\text{CO}_2\text{H}$	644.2 (M + H)	3.63

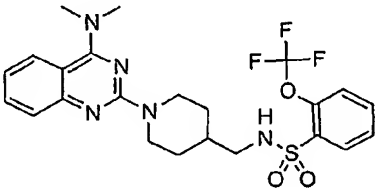
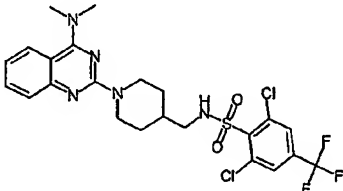
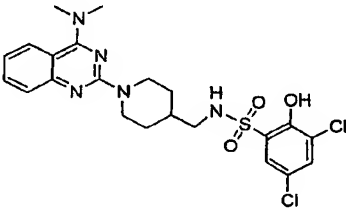
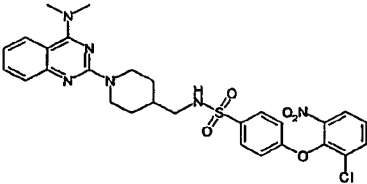
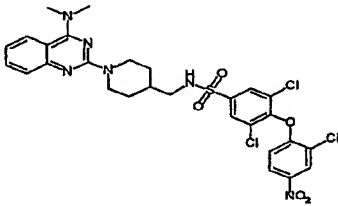
Example No.	Structure	ESI-MS	Retention Time (min)
3545	 $\text{CF}_3\text{CO}_2\text{H}$	520.4 (M + H)	3.56
3546	 $\text{CF}_3\text{CO}_2\text{H}$	504.2 (M + H)	3.25
3547	 $2\text{CF}_3\text{CO}_2\text{H}$	513.4 (M + H)	2.86
3548	 $\text{CF}_3\text{CO}_2\text{H}$	616.2 (M + H)	3.73
3549	 $2\text{CF}_3\text{CO}_2\text{H}$	450.4 (M + H)	2.79
3550	 $\text{CF}_3\text{CO}_2\text{H}$	466.2 (M + H)	3.35

Example No.	Structure	ESI-MS	Retention Time (min)
3551	 $2\text{CF}_3\text{CO}_2\text{H}$	465.2 (M + H)	3.34
3552	 $\text{CF}_3\text{CO}_2\text{H}$	451.2 (M + H)	3.83
3553	 $\text{CF}_3\text{CO}_2\text{H}$	451.2 (M + H)	4.10
3554	 $\text{CF}_3\text{CO}_2\text{H}$	563.2 (M + H)	4.33
3555	 $2\text{CF}_3\text{CO}_2\text{H}$	468.4 (M + H)	3.66
3556	 $2\text{CF}_3\text{CO}_2\text{H}$	467.4 (M + H)	2.85

Example No.	Structure	ESI-MS	Retention Time (min)
3557	 $\text{CF}_3\text{CO}_2\text{H}$	515.4 (M + H)	3.52
3558	 $\text{CF}_3\text{CO}_2\text{H}$	485.2 (M + H)	3.40
3559	 $2\text{CF}_3\text{CO}_2\text{H}$	467.4 (M + H)	3.90
3560	 $\text{CF}_3\text{CO}_2\text{H}$	473.4 (M + H)	4.17
3561	 $\text{CF}_3\text{CO}_2\text{H}$	467.4 (M + H)	3.57
3562	 $\text{CF}_3\text{CO}_2\text{H}$	490.2 (M + H)	4.00

Example No.	Structure	ESI-MS	Retention Time (min)
3563	 <chem>COc1ccc(cc1)C(=O)Nc2ccc(Nc3nc4ccccc4n3C)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	490.2 (M + H)	3.99
3564	 <chem>c1ccc(cc1)c2nc3ccccc3nc2C(=O)Nc4ccc(Nc5nc6ccccc6n5C)cc4</chem> $2\text{CF}_3\text{CO}_2\text{H}$	476.2 (M + H)	3.76
3565	 <chem>c1ccc(cc1)c2sc(C(=O)Nc3ccc(Nc4nc5ccccc5n4C)cc3)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	467.2 (M + H)	4.07
3566	 <chem>Clc1cc(Cl)ccc1C(=O)Nc2ccc(Nc3nc4ccccc4n3C)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	528.2 (M + H)	4.53
3567	 <chem>c1ccc(cc1)c2oc(C(=O)Nc3ccc(Nc4nc5ccccc5n4C)cc3)cc2</chem> $\text{CF}_3\text{CO}_2\text{H}$	464.2 (M + H)	4.11
3568	 <chem>Clc1cc(Cl)ccc1S(=O)(=O)NC2CCCN2c3nc4ccccc4n3C</chem> $\text{CF}_3\text{CO}_2\text{H}$	494.0 (M + H)	3.43

Example No.	Structure	ESI-MS	Retention Time (min)
3569	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4ccc(F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	444.0 (M + H)	3.03
3570	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4ccc(I)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	552.0 (M + H)	3.30
3571	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4ccc(OC(F)(F)F)cc4</chem> $\text{CF}_3\text{CO}_2\text{H}$	510.0 (M + H)	3.37
3572	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4ccc(C(F)(F)F)cc4C(F)(F)F</chem> $\text{CF}_3\text{CO}_2\text{H}$	562.0 (M + H)	3.66
3573	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4cc(OC(F)(F)F)cc(OC(F)(F)F)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	622.0 (M + H)	3.61
3574	 <chem>CC1=NC2=CC=CC=C2N(C)=N1CCN3CCCCC3CCNS(=O)(=O)c4cc(OC(F)(F)F)cc(Br)c4</chem> $\text{CF}_3\text{CO}_2\text{H}$	588.0 (M + H)	3.59

Example No.	Structure	ESI-MS	Retention Time (min)
3575	 <chem>CC1=NC2=CC=CC=C2N1N=C3C(=NC(=CC=C3)N(C)C)S(=O)(=O)NCC4CCNCC4c5ccccc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	510.0 (M + H)	3.31
3576	 <chem>CC1=NC2=CC=CC=C2N1N=C3C(=NC(=CC=C3)N(C)C)S(=O)(=O)NCC4CCNCC4c5cc(Cl)c(Cl)c(C(F)(F)F)c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	562.0 (M + H)	3.61
3577	 <chem>CC1=NC2=CC=CC=C2N1N=C3C(=NC(=CC=C3)N(C)C)S(=O)(=O)NCC4CCNCC4c5cc(Cl)c(Cl)cc5O</chem> $\text{CF}_3\text{CO}_2\text{H}$	510.0 (M + H)	3.35
3578	 <chem>CC1=NC2=CC=CC=C2N1N=C3C(=NC(=CC=C3)N(C)C)S(=O)(=O)NCC4CCNCC4c5ccc(Cl)c([N+](=O)[O-])c5</chem> $\text{CF}_3\text{CO}_2\text{H}$	597.0 (M + H)	3.55
3579	 <chem>CC1=NC2=CC=CC=C2N1N=C3C(=NC(=CC=C3)N(C)C)S(=O)(=O)NCC4CCNCC4c5cc(Cl)c(Oc6cc(Cl)c(Cl)cc6)cc5</chem> $\text{CF}_3\text{CO}_2\text{H}$	665.0 (M + H)	4.02

Assay Procedures

Compounds identified and disclosed throughout this patent document were assayed according to the protocols found in co-pending patent application having U.S. Serial Number 09/826,509, which is incorporated herein by reference.

Example 3580

Preparation of Endogenous MCH Receptor.

The endogenous human MCH receptor was obtained by PCR using genomic DNA as template and rTth polymerase (Perkin Elmer) with the buffer system provided by the manufacturer, 0.25 μ M of each primer, and 0.2 mM of each 4 nucleotides. The cycle condition was 30 cycles of 94°C for 1 min, 56°C for 1min and 72 °C for 1 min and 20 sec. The 5' PCR primer contained a HindIII site with the sequence:

5'-GTGAAGCTTGCCTCTGGTGCCTGCAGGAGG-3' (SEQ.ID.NO.:1)

and the 3' primer contained an EcoRI site with the sequence:

5'-GCAGAATCCCGGTGGCGTGTTGTGGTGCCC-3' (SEQ.ID.NO.:2).

The 1.3 kb PCR fragment was digested with HindIII and EcoRI and cloned into HindIII-EcoRI site of CMVp expression vector. Later the cloning work by Lakaye et al showed that there is an intron the coding rgion of the gene. Thus the 5' end of the cDNA was obtained by 5' RACE PCR using Clontech's marathon-ready hypothalamus cDNA as template and the manufacturer's recommended protocol for cycling condition. The 5' RACE PCR for the first and second round PCR were as follows:

5'-CATGAGCTGGTGGATCATGAAGGG-3' (SEQ.ID.NO.:3) and

5'-ATGAAGGGCATGCCAGGAGAAAG-3' (SEQ.ID.NO.:4).

Nucleic acid and amino acid sequences were thereafter determined and verified with the published sequences found on GenBank having Accession Number U71092.

Example 3581

Preparation of Non-Endogenous, Constitutively Active MCH Receptor.

Preparation of a non-endogenous version of the human MCH receptor was accomplished by creating a MCH-IC3-SST2 mutation (*see*; SEQ.ID.NO.:7 for nucleic acid sequence, and SEQ.ID.NO.:8 for amino acid sequence). Blast result showed that MCH receptor had the highest sequence homology to known SST2 receptor. Thus the third intracellular loop ("IC3") of MCH receptor was replaced with that of the IC3 of SST2

receptor to see if the chimera would show constitutive activity.

The BamHI-BstEII fragment containing IC3 of MCH receptor was replaced with synthetic oligonucleotides that contained the IC3 of SST2. The PCR sense mutagenesis primer used had the following sequence:

5'-GATCCTGCAGAAGGTGAAGTCCTCTGGAATCCGAGTGGGCTCCTCTAAGAG
GAAGAAGTCTGAGAAGAAG-3' (SEQ.ID.NO.:9)

and the antisense primer had the following sequence:

5'-GTGACCTTCTTCTCAGACTTCTTCCTCTTAGAGGAGCCCACTCGGATTCCAG
AGGACTTCACCTTCTGCAG-3' (SEQ.ID.NO.:10).

The endogenous MCH receptor cDNA was used as a template.

Example 3582

GPCR Fusion Protein Preparation.

MCH Receptor-Gi α Fusion Protein construct was made as follows: primers were designed for endogenous MCH receptor was as follows:

5'-GTGAAGCTTGCCCGGCAGGATGGACCTGG-3' (SEQ.ID.NO.:11; sense)

5'-ATCTAGAGGTGCCTTTGCTTTCTG-3' (SEQ.ID.NO.:12; antisense).

The sense and anti-sense primers included the restriction sites for KB4 and XbaI, respectively.

PCR was utilized to secure the respective receptor sequences for fusion within the Gi α universal vector disclosed above, using the following protocol for each: 100ng cDNA for MCH receptor was added to separate tubes containing 2ul of each primer (sense and anti-sense), 3uL of 10mM dNTPs, 10uL of 10XTaqPlus™ Precision buffer, 1uL of TaqPlus™ Precision polymerase (Stratagene: #600211), and 80uL of water. Reaction temperatures and cycle times for MCH receptor were as follows: the initial denaturing step was done at 94°C for five minutes, and a cycle of 94°C for 30 seconds; 55°C for 30 seconds; 72°C for two minutes. A final extension time was done at 72°C for ten minutes. PCR product was run on a 1% agarose gel and then purified (data not shown). The purified product was digested with KB4 and XbaI (New England Biolabs) and the desired inserts will be isolated, purified and ligated into the Gi universal vector at the respective restriction site. The positive clones were isolated following transformation and determined by restriction enzyme digest; expression using 293 cells was accomplished.

following the protocol set forth *infra*. Each positive clone for MCH receptor: Gi-Fusion Protein was sequenced and made available for the direct identification of candidate compounds. (See, SEQ.ID.NO.:13 for nucleic acid sequence and SEQ.ID.NO.:14 for amino acid sequence).

Endogenous version of MCH receptor was fused upstream from the G protein Gi and is located at nucleotide 1 through 1,059 (see, SEQ.ID.NO.:13) and amino acid residue 1 through 353 (see, SEQ.ID.NO.:14). With respect to the MCH receptor, 2 amino acid residues (an equivalent of 6 nucleotides) were placed in between the endogenous (or non-endogenous) GPCR and the start codon for the G protein $G_{i\alpha}$. Therefore, the Gi protein is located at nucleotide 1,066 through 2,133 (see, SEQ.ID.NO.:13) and at amino acid residue 356 through 711 (see, SEQ.ID.NO.:14). Those skilled in the art are credited with the ability to select techniques for constructing a GPCR Fusion Protein where the G protein is fused to the 3' end of the GPCR of interest.

Example 3583

ASSAY FOR DETERMINATION OF CONSTITUTIVE ACTIVITY OF NON-ENDOGENOUS GPCRS

A. Intracellular IP_3 Accumulation Assay

On day 1, cells comprising the receptors (endogenous and/or non-endogenous) can be plated onto 24 well plates, usually 1×10^5 cells/well (although this number can be optimized). On day 2 cells can be transfected by firstly mixing 0.25ug DNA in 50 ul serum free DMEM/well and 2 ul lipofectamine in 50 ul serum-free DMEM/well. The solutions are gently mixed and incubated for 15-30 min at room temperature. Cells are washed with 0.5 ml PBS and 400 ul of serum free media is mixed with the transfection media and added to the cells. The cells are then incubated for 3-4 hrs at $37^\circ\text{C}/5\%\text{CO}_2$ and then the transfection media is removed and replaced with 1ml/well of regular growth media. On day 3 the cells are labeled with ^3H -myo-inositol. Briefly, the media is removed and the cells are washed with 0.5 ml PBS. Then 0.5 ml inositol-free/serum free media (GIBCO BRL) is added/well with 0.25 μCi of ^3H -myo-inositol/ well and the cells are incubated for 16-18 hrs o/n at $37^\circ\text{C}/5\%\text{CO}_2$. On Day 4 the cells are washed with 0.5 ml PBS and 0.45 ml of assay medium is added containing inositol-free/serum free media 10 μM pargyline 10 mM lithium chloride or 0.4 ml of assay medium and 50 ul of 10x

ketanserin (ket) to final concentration of 10 μ M. The cells are then incubated for 30 min at 37°C. The cells are then washed with 0.5 ml PBS and 200 μ l of fresh/ice cold stop solution (1M KOH; 18 mM Na-borate; 3.8 mM EDTA) is added/well. The solution is kept on ice for 5-10 min or until cells were lysed and then neutralized by 200 μ l of fresh/ice cold neutralization sol. (7.5 % HCL). The lysate is then transferred into 1.5 ml eppendorf tubes and 1 ml of chloroform/methanol (1:2) is added/tube. The solution is vortexed for 15 sec and the upper phase is applied to a Biorad AG1-X8™ anion exchange resin (100-200 mesh). Firstly, the resin is washed with water at 1:1.25 W/V and 0.9 ml of upper phase is loaded onto the column. The column is washed with 10 mls of 5 mM myo-inositol and 10 ml of 5 mM Na-borate/60mM Na-formate. The inositol tris phosphates are eluted into scintillation vials containing 10 ml of scintillation cocktail with 2 ml of 0.1 M formic acid/ 1 M ammonium formate. The columns are regenerated by washing with 10 ml of 0.1 M formic acid/3M ammonium formate and rinsed twice with H₂O and stored at 4°C in water.

Reference is made to Figure 1. Figure 1 provides an illustration of IP₃ production from several non-endogenous, constitutively activated version of MCH receptor as compared with the endogenous version of this receptor. When compared to the endogenous version of MCH receptor ("MCH-R wt"), MCH-IC3-SST2 evidenced about a 27% increase in IP₃ accumulation.

Example 3584

Determination of Compound Using [³⁵S]GTP γ S ASSAY

Direct identification of candidate compounds was initially screened using [³⁵S]GTP γ S Assay (see, Example 6 of co-pending patent application 09/826,509). Preferably, an MCH receptor: Gi Fusion Protein was utilized, according to Example 6(2) of co-pending patent application 09/826,509. Several lead hits were identified utilizing [³⁵S]GTP γ S Assay.

Example 3585

High Throughput Functional Screening: FLIPR™

Subsequently, a functional based assay was used to confirm the lead hits, referred to as FLIPR™ (the Fluorometric Imaging Plate Reader) and FDSS6000™ (Functional

Drug Screening System). This assay utilized a non-endogenous version of the MCH receptor, which was created by swapping the third intracellular loop of the MCH receptor with that of the SST2 receptor (see Example 2(B)(2) of patent application serial number 09/826,509).

The FLIPR and FDSS assays are able to detect intracellular Ca^{2+} concentration in cells, which can be utilized to assess receptor activation and determine whether a candidate compound is an, for example, antagonist, inverse agonist or agonist to a Gq-coupled receptor. The concentration of free Ca^{2+} in the cytosol of any cell is extremely low, whereas its concentration in the extracellular fluid and endoplasmic reticulum (ER) is very high. Thus, there is a large gradient tending to drive Ca^{2+} into the cytosol across both the plasma membrane and ER. The FLIPRTM and FDSS6000TM systems (Molecular Devices Corporation, HAMAMATSU Photonics K.K.) are designed to perform functional cell-based assays, such as the measurement of intracellular calcium for high-throughput screening. The measurement of fluorescent is associated with calcium release upon activation of the Gq-coupled receptors. Gi or Go coupled receptors are not as easily monitored through the FLIPRTM and FDSS6000TM systems because these G proteins do not couple with calcium signal pathways.

To confirm the lead hits identified using the [³⁵S]GTP γ S assay, Fluorometric Imaging Plate Reader system was used to allow for rapid, kinetic measurements of intracellular fluorescence in 96 well microplates (or 384 well microplates). Simultaneous measurements of fluorescence in all wells can be made by FLIPR or FDSS6000TM every second with high sensitivity and precision. These systems are ideal for measuring cell-based functional assays such as monitoring the intracellular calcium fluxes that occur within seconds after activation of the Gq coupled receptor.

Briefly, the cells are seeded into 96 well at 5.5×10^4 cells/well with complete culture media (Dulbecco's Modified Eagle Medium with 10 % fetal bovine serum, 2 mM L-glutamine, 1 mM sodium pyruvate and 0.5 mg/ml G418, pH 7.4) for the assay next day. On the day of assay, the media is removed and the cells are incubated with 100 μ l of loading buffer (4 μ M Fluo4-AM in complete culture media containing 2.5 mM Probenicid, 0.5 mg/ml and 0.2% bovine serum albumin) in 5% CO₂ incubator at 37°C for 1 hr. The loading buffer is removed, and the cells are washed with wash buffer (Hank's Balanced Salt Solution containing 2.5 mM Probenicid, 20 mM HEPES, 0.5 mg/ml and 0.2% bovine

serum albumin, pH 7.4)). One hundred fifty μ l of wash buffer containing various concentrations of test compound are added to the cells, and the cells are incubated in 5% CO₂ incubator at 37°C for 30 min. Fifty μ l of wash buffer containing various concentration of MCH are added to each well, and transient changes in [Ca²⁺]_i evoked by MCH are monitored using the FLIPR or FDSS in 96 well plates at Ex. 488 nm and Em. 530 nm for 290 second. When antagonist activity of compound is tested, 50 nM of MCH is used.

Use of FLIPR™ and FDSS6000™ can be accomplished by following manufacturer's instruction (Molecular Device Corporation and HAMAMATSU Photonics K.K.).

The results were shpwn below.

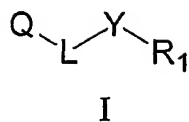
Compound No.	IC ₅₀ value (nM)
Example 41	6
Example 42	19

It is intended that each of the patents, applications, printed publications, and other published documents mentioned or referred to in this specification be herein incorporated by reference in their entirety.

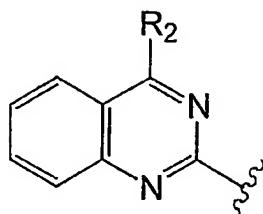
Those skilled in the art will appreciate that numerous changes and modifications may be made to the preferred embodiments of the invention and that such changes and modifications may be made without departing from the spirit of the invention. It is therefore intended that the appended claims cover all such equivalent variations as fall within the true spirit and scope of the invention.

What is claimed is:

1. A compound of Formula I:

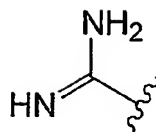


wherein Q is



II

or



III

R₁ represents

(i) C₁-C₁₆ alkyl,

C₁-C₁₆ alkyl substituted by substituent(s) independently selected from

- halogen,
- hydroxy,
- oxo,
- C₁-C₃ alkoxy,
- C₁-C₃ alkoxy substituted by substituent(s) independently selected from
- carbocyclic aryl,
- heterocyclyl,
- heterocyclyl substituted by C₁-C₃ alkyl,
- C₁-C₃ alkylcarbonyloxy,
- carbocyclyloxy,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by substituent(s) independently selected from
- halogen,
- nitro,
- carbocyclic aryl,
- carbocyclic aryl substituted by C₁-C₃ alkoxy,

- C₁-C₄ alkyl,
- C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - oxo,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,
 - mono- or di-C₁-C₃ alkylamino substituted by halogenated carbocyclic aryl,
 - carbocyclic arylcarbonylamino,
 - halogenated carbocyclic arylcarbonylamino,
 - heterocyclyloxy,
 - heterocyclyloxy substituted by C₁-C₃ alkyl,
 - substituted heterocyclyl-ethylideneaminooxy,
 - C₁-C₃ alkoxy carbonyl,
 - C₁-C₃ alkoxy carbonyl substituted by carbocyclic aryl,
 - mono- or di-C₁-C₃ alkylaminocarbonyl,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-C₁-C₃ alkylamino substituted by substituent(s) independently selected from
 - cyano,
 - carbocyclic aryl,
 - heterocyclyl,
 - mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylamino substituted by substituent(s) independently selected from
 - hydroxy,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkylcarbonylamino,
 - C₁-C₃ alkylcarbonylamino substituted by substituent(s) independently selected from
 - C₁-C₃ alkylcarbonylamino,
 - carbocyclic arylcarbonylamino,
 - heterocyclyl,
 - C₁-C₄ alkoxy carbonylamino,
 - heterocyclyl carbonylamino,
 - carbocyclic arylsulfonylamino,

- carbocyclic arylsulfonylamino substituted by substituent(s) independently selected from
 - nitro,
 - C₁-C₃ alkyl,
 - mono- or di-C₁-C₃ alkylamino,
- C₁-C₃ alkylthio,
- C₁-C₃ alkylthio substituted by substituent(s) independently selected from
 - mono- or di-carbocyclic arylaminocarbonyl,
 - halogenated mono- or di-carbocyclic arylaminocarbonyl,
 - mono- or di-carbocyclic arylamino,
 - halogenated mono- or di-carbocyclic arylamino,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkoxy,
- carbocyclic arylthio,
- carbocyclic arylthio substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
- carbocyclic arylsulfonyl,
- halogenated carbocyclic arylsulfonyl,
- heterocyclylthio,
- heterocyclylthio substituted by substituent(s) independently selected from
 - nitro,
 - C₁-C₃ alkyl,
 - C₃-C₆ cycloalkyl,
 - C₃-C₆ cycloalkyl substituted by C₁-C₃ alkyl,
 - C₃-C₆ cycloalkenyl,
- carbocyclyl,
- carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,

- C₂-C₃ alkenyl,
- C₂-C₃ alkenyl substituted by carbocyclic aryl,
- C₂-C₃ alkenyl substituted by carbocyclic aryl substituted C₁-C₃ alkylsulfinyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - nitro,
- C₁-C₄ alkyl,
- C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - oxo,
 - carbocyclic aryl,
 - heterocyclyl,
 - mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylamino substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
- C₁-C₄ alkoxy,
- C₁-C₄ alkoxy substituted by substituent(s) independently selected from
 - halogen,
 - carbocyclic aryl,
 - carbocyclic aryloxy,
 - C₁-C₃ alkoxycarbonyl,
 - C₁-C₃ alkylcarbonyloxy,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-carbocyclic arylamino,

- halogenated mono- or di-carbocyclic arylamino,
- mono- or di-carbocyclic arylaminocarbonyl,
- mono- or di-carbocyclic arylaminocarbonyl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
 - mercapto,
 - C₁-C₃ alkylthio,
 - halogenated C₁-C₃ alkylthio,
 - C₁-C₃ alkylsulfonyl,
 - C₃-C₆ cycloalkyl,
 - carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - hydroxy,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkyl substituted by carbocyclic aryl,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (ii) C₂-C₈ alkenyl,
C₂-C₈ alkenyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,

•carbocyclic aryl substituted by substituent(s) independently selected from

••halogen,

••hydroxy,

••nitro,

••C₁-C₃ alkyl,

••halogenated C₁-C₃ alkyl,

••C₁-C₃ alkoxy,

••halogenated C₁-C₃ alkoxy,

•heterocyclyl,

•heterocyclyl substituted by substituent(s) independently selected from

••hydroxy,

••nitro,

••C₁-C₃ alkyl,

••C₁-C₃ alkoxy,

(iii) C₂-C₄ alkynyl,

C₂-C₄ alkynyl substituted by carbocyclic aryl,

(iv) C₃-C₆ cycloalkyl,

C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from

•C₁-C₃ alkyl,

•C₁-C₃ alkyl substituted by substituent(s) independently selected from

••hydroxy,

••oxo,

••carbocyclic aryl,

•mono- or di-C₁-C₃ alkylamino,

•mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,

•carbocyclic arylcarbonylamino,

•carbocyclic aryl,

(v) C₃-C₆ cycloalkyl,

C₃-C₆ cycloalkyl substituted by C₁-C₃ alkyl,

(vi) carbocyclyl,

carbocyclyl substituted by substituent(s) independently selected from

•hydroxy,

- nitro,
- (vii) carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₉ alkyl,
 - C₁-C₉ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - oxo,
 - C₁-C₃ alkoxy,
 - carbocyclic aryloxy,
 - mono- or di-C₁-C₃ alkylamino-N-oxy,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,
 - mono- or di-carbocyclic arylamino,
 - carbocyclylimino,
 - carbocyclylimino substituted by carbocyclic aryl,
 - mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylamino substituted by C₁-C₃ alkoxy,
 - mono- or di-carbocyclic arylaminocarbonyl,
 - mono- or di-carbocyclic arylaminocarbonyl substituted by C₁-C₃ alkoxy,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,
 - heterocyclyl,
 - heterocyclyl substituted by C₁-C₃ alkyl,
 - C₂-C₃ alkenyl,

- C₂-C₃ alkenyl substituted by carbocyclic aryl,
- C₁-C₉ alkoxy,
- C₁-C₉ alkoxy substituted by substituent(s) independently selected from
 - hydroxy,
 - halogen,
 - carboxy,
 - mono- or di-C₁-C₃ alkylamino,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,
- C₂-C₃ alkenyloxy,
- C₁-C₃ alkylcarbonyloxy,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,
 - halogenated C₁-C₄ alkyl,
 - C₁-C₃ alkoxy,
 - heterocyclyloxy,
 - heterocyclyloxy substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,
- (carbocyclic aryl)S(O)₂O,

- carboxy,
- C₁-C₃ alkoxycarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl substituted by carbocyclic aryl,
- mono- or di-carbocyclic arylaminocarbonyl,
- mono- or di-carbocyclic arylaminocarbonyl substituted by C₁-C₃ alkyl,
- amino,
- mono- or di-C₁-C₄ alkylamino,
- mono- or di-C₁-C₄ alkylamino substituted by cyano,
- mono- or di-carbocyclic arylamino,
- C₁-C₃ alkynylcarbonylamino,
- C₁-C₃ alkynylcarbonylamino substituted by carbocyclic aryl,
- carbocyclic arylsulfonylamino,
- carbocyclic arylsulfonylamino substituted by C₁-C₃ alkyl,
- (carbocyclic aryl)NHC(O)NH,
- (carbocyclic aryl)NHC(O)NH substituted by C₁-C₃ alkoxy,
- (carbocyclic aryl)NHC(O)NH substituted by halogenated C₁-C₃ alkoxy,
- carbocyclic aryl diazo,
- carbocyclic aryl diazo substituted by mono- or di- C₁-C₃ alkylamino,
- C₁-C₃ alkylthio,
- halogenated C₁-C₃ alkylthio,
- carbocyclic arylthio,
- carbocyclic arylthio substituted by substituent(s) independently selected from
 - halogen,
 - cyano,
 - C₁-C₃ alkyl,
 - heterocyclylthio,
 - C₁-C₃ alkylsulfonyl,
 - mono- or di-C₁-C₃ alkylaminosulfonyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - C₁-C₇ alkyl,

- halogenated C₁-C₇ alkyl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (viii) heterocyclyl,
or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - oxo,
 - C₁-C₃ alkylcarbonyloxy,
 - carbocyclic arylcarbonylamino,
 - halogenated carbocyclic arylcarbonylamino,
 - C₁-C₃ alkoxycarbonyl,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkylthio substituted by carbocyclic aryl,
 - C₁-C₃ alkylthio substituted by halogenated carbocyclic aryl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,

- halogenated C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- C₁-C₃ alkoxy substituted by carbocyclic aryl,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - mono- or di-C₁-C₃ alkylamino,
 - C₁-C₄ alkylcarbonylamino,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkenylthio,
 - carbocyclic arylthio,
 - halogenated carbocyclic arylthio,
 - carbocyclic arylthio substituted by C₁-C₃ alkoxycarbonyl,
 - heterocyclylthio,
 - heterocyclylthio substituted by C₁-C₃ alkyl,
 - C₁-C₃ alkylsulfonyl,
 - carbocyclic arylsulfonyl,
 - halogenated carbocyclic arylsulfonyl,
 - carbocyclic arylsulfonyl substituted by C₁-C₄ alkyl,
 - C₁-C₃ alkoxycarbonyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - halogen,

- C₁-C₃ alkyl,
- halogenated C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- C₁-C₃ alkoxycarbonyl;

R₂ is -NHNH₂, -NHNHBoc, -N(R_{2a})(R_{2b}), morpholino, 4-acetyl-piperazyl, or 4-phenyl-piperazyl;

wherein R_{2a} is H or C₁-C₃ alkyl;

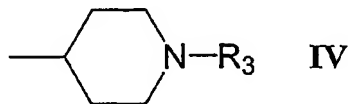
R_{2b} is C₁-C₄ alkyl, C₁-C₄ alkyl substituted by substituent(s) independently selected from

- hydroxy,
- C₁-C₃ alkoxy,
- amino,
- NHBoc,
- C₃-C₆ cycloalkyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
- halogen,
- C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- SO₂NH₂,
- heterocyclyl,

C₃-C₆ cycloalkyl, carbocyclic aryl, carbocyclic aryl substituted by substituent(s) independently selected from

- halogen,
- C₁-C₃ alkyl,
- C₁-C₃ alkoxy,

or a group of Formula IV;

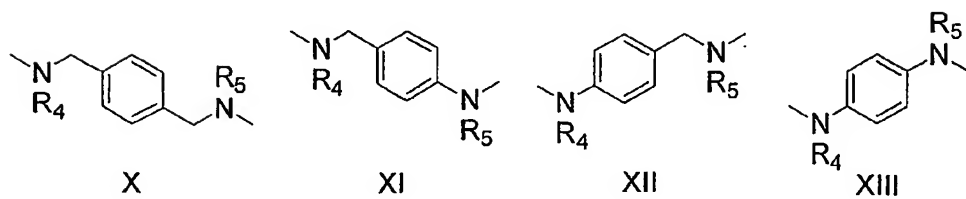
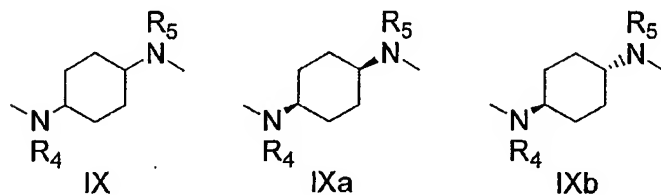
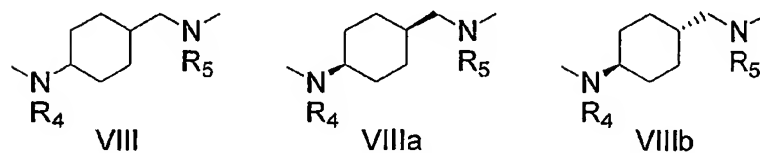
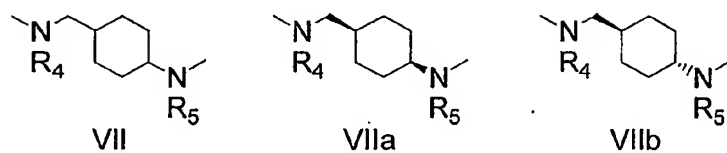
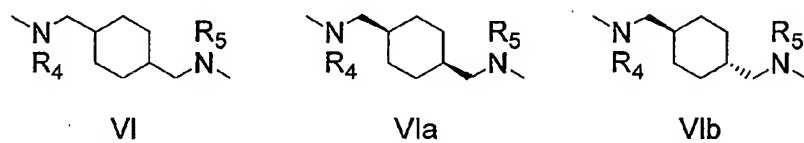
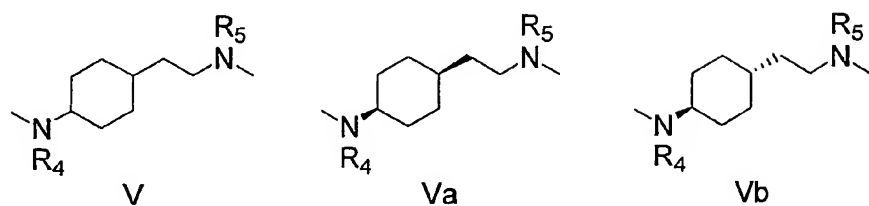


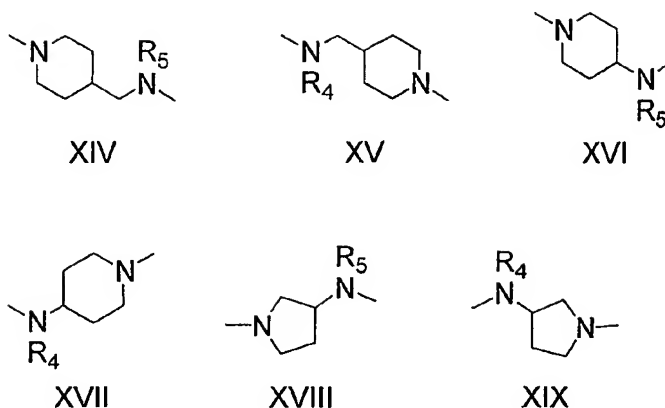
wherein Boc is carbamic acid *tert*-butyl ester and R₃ is C₁-C₃ alkyl or C₁-C₃ alkyl

substituted by substituent(s) independently selected from

- carbocyclic aryl,
- halogenated carbocyclic aryl,
- carbocyclic aryl substituted by C₁-C₃ alkoxy;

L is selected from Formula V - XIX;





wherein R_4 is H or C_1 - C_3 alkyl;

R_5 is H, C_1 - C_3 alkyl, or C_1 - C_3 alkyl substituted by a substituted carbocyclic aryl;

Y is $-S(O)_2-$, $-C(O)-$, or $-(CH_2)_m$;

m is 0 or 1;

wherein carbocyclic aryl is phenyl, naphthyl, anthranyl, biphenyl, or phenanthryl;

carbocyclyl is 10,11-dihydro-5-oxo-dibenzo[a,d]cycloheptyl, 1-oxo-indanyl, 7,7-dimethyl-2-oxo-bicyclo[2.2.1]heptyl, 9H-fluorenyl, 9-oxo-fluorenyl, acenaphthyl, anthraquinonyl, C-fluoren-9-ylidene, indanyl, indenyl, 1,2,3,4-tetrahydro-naphthyl, or bicyclo[2.2.1]hepteny;

heterocyclyl is 1,2,3,4-tetrahydro-isoquinolyl, 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1,3,4-thiadiazolyl, 1,3-dioxo-isoindolyl, 1,3-dioxolanyl, 1H-indolyl, 1H-pyrrolo[2,3-c]pyridyl, 1H-pyrrolyl, 1-oxo-3H-isobenzofuranyl, 2,2',5',2"-terthiophenyl, 2,2'-bithiophenyl, 2,3-dihydro-1-oxo-isoindolyl, 2,3-dihydro-benzo[1,4]dioxinyl, 2,4-dihydro-3-oxo-pyrazolyl, 2H-benzopyranyl, 2-oxo-benzopyranyl, 2-oxo-pyrrolidinyl, 3,4-dihydro-2H-benzo[1,4]oxazinyl, 3,4-dihydro-2H-benzo[b][1,4]dioxepinyl, 4H-benzo[1,3]dioxinyl, 4H-benzopyranyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, 4-oxo-3,4-dihydro-phthalazinyl, 4-oxo-benzopyranyl, 9,10,10-trioxo-thioxanthenyl, 9H-carbazolyl, 9H-xanthenyl, azetidiny, benzimidazolyl, benzo[1,3]dioxolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, benzofuryl, benzothiazolyl, cinnolyl, furyl, imidazo[2,1-b]thiazolyl, imidazolyl, isoxazolyl, morpholino, morpholinyl, oxazolyl, oxolanyl, piperazyl, piperidyl, piridyl, pyrazolo[5,1-b]thiazolyl, pyrazolyl, pyridyl, pyrimidyl, pyrrolidyl, quinolyl, quinoxalyl, thiazolidyl, thiazolyl, thienyl, thiolanyl, 2,3-

dihydro-benzofuryl, tetrahydro-thienyl, or benzofuranyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

2. A compound according to claim 1, wherein Q is Formula II;

R₁ represents

(i) C₁-C₁₀ alkyl,

C₁-C₁₀ alkyl substituted by substituent(s) independently selected from

•halogen,

•oxo,

•C₁-C₃ alkoxy,

•C₁-C₃ alkoxy substituted by carbocyclic aryl,

•C₁-C₃ alkylcarbonyloxy,

•carbocyclyloxy,

•carbocyclic aryloxy,

•carbocyclic aryloxy substituted by substituent(s) independently selected from

••halogen,

••nitro,

••C₁-C₄ alkyl,

••C₁-C₄ alkyl substituted by substituent(s) independently selected from

•••oxo,

•••carbocyclic arylcarbonylamino,

•••halogenated carbocyclic arylcarbonylamino,

•heterocyclyloxy,

•heterocyclyloxy substituted by C₁-C₃ alkyl,

•substituted heterocyclyl-ethylideneaminooxy,

•C₁-C₃ alkoxycarbonyl,

•C₁-C₃ alkoxycarbonyl substituted by carbocyclic aryl,

•mono- or di-C₁-C₃ alkylaminocarbonyl,

•mono- or di-carbocyclic arylamino,

•mono- or di-carbocyclic arylamino substituted by hydroxy,

•C₁-C₃ alkylcarbonylamino,

- C₁-C₃ alkylcarbonylamino substituted by substituent(s) independently selected from
 - C₁-C₃ alkylcarbonylamino,
 - carbocyclic arylcarbonylamino,
 - heterocyclyl,
- C₁-C₄ alkoxycarbonylamino,
- heterocyclyl carbonylamino,
- carbocyclic arylsulfonylamino,
- carbocyclic arylsulfonylamino substituted by substituent(s) independently selected from
 - nitro,
 - C₁-C₃ alkyl,
 - mono- or di-C₁-C₃ alkylamino,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkylthio substituted by substituent(s) independently selected from
 - mono- or di-carbocyclic arylaminocarbonyl,
 - halogenated mono- or di-carbocyclic arylaminocarbonyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkoxy,
 - carbocyclic arylthio,
 - carbocyclic arylthio substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - carbocyclic arylsulfonyl,
 - halogenated carbocyclic arylsulfonyl,
 - heterocyclylthio,
 - heterocyclylthio substituted by substituent(s) independently selected from
 - nitro,
 - C₁-C₃ alkyl,
 - C₃-C₆ cycloalkyl,
 - C₃-C₆ cycloalkyl substituted by C₁-C₃ alkyl,
 - C₃-C₆ cycloalkenyl,

- carbocyclyl,
- carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - C₂-C₃ alkenyl,
 - C₂-C₃ alkenyl substituted by carbocyclic aryl,
 - C₂-C₃ alkenyl substituted by carbocyclic aryl substituted C₁-C₃ alkylsulfinyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - oxo,
 - carbocyclic aryl,
 - heterocyclyl,
 - C₁-C₄ alkoxy,
 - C₁-C₄ alkoxy substituted by substituent(s) independently selected from
 - halogen,
 - carbocyclic aryl,
 - carbocyclic aryloxy,
 - C₁-C₃ alkylcarbonyloxy,
 - mono- or di-carbocyclic arylamino,
 - halogenated mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylaminocarbonyl,
 - mono- or di-carbocyclic arylaminocarbonyl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,

- C₁-C₃ alkoxy,
- halogenated C₁-C₃ alkoxy,
- mercapto,
- C₁-C₃ alkylthio,
- halogenated C₁-C₃ alkylthio,
- C₁-C₃ alkylsulfonyl,
- C₃-C₆ cycloalkyl,
- carbocyclic aryl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
 - hydroxy,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkyl substituted by carbocyclic aryl,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (ii) C₂-C₆ alkenyl,
 - C₂-C₆ alkenyl substituted by substituent(s) independently selected from
 - oxo,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - hydroxy,

- C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- (iii) C₃-C₆ cycloalkyl,
C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - C₁-C₃ alkyl substituted by substituent(s) independently selected from
 - oxo,
 - carbocyclic aryl,
 - carbocyclic arylcarbonylamino,
 - carbocyclic aryl,
 - (iv) carbocyclyl,
carbocyclyl substituted by nitro,
 - (v) carbocyclic aryl,
carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₉ alkyl,
 - C₁-C₉ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - carbocyclic aryloxy,
 - carbocyclylimino,
 - carbocyclylimino substituted by carbocyclic aryl,
 - mono- or di-carbocyclic arylaminocarbonyl,
 - mono- or di-carbocyclic arylaminocarbonyl substituted by C₁-C₃ alkoxy,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl,

- heterocyclyl,
- heterocyclyl substituted by C₁-C₃ alkyl,
- C₁-C₇ alkoxy,
- C₁-C₇ alkoxy substituted by substituent(s) independently selected from
- halogen,
- carbocyclic aryl,
- C₁-C₃ alkylcarbonyloxy,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by C₁-C₃ alkoxy,
- C₁-C₃ alkoxycarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl substituted by carbocyclic aryl,
- mono- or di-carbocyclic arylaminocarbonyl,
- mono- or di-carbocyclic arylaminocarbonyl substituted by C₁-C₃ alkyl,
- amino,
- mono- or di-C₁-C₃ alkylamino,
- C₁-C₃ alkynylcarbonylamino,
- C₁-C₃ alkynylcarbonylamino substituted by carbocyclic aryl,
- carbocyclic arylsulfonylamino,
- carbocyclic arylsulfonylamino substituted by C₁-C₃ alkyl,
- (carbocyclic aryl)NHC(O)NH,
- (carbocyclic aryl)NHC(O)NH substituted by C₁-C₃ alkoxy,
- (carbocyclic aryl)NHC(O)NH substituted by halogenated C₁-C₃ alkoxy,
- C₁-C₃ alkylthio,
- halogenated C₁-C₃ alkylthio,
- carbocyclic arylthio,
- carbocyclic arylthio substituted by cyano,
- C₁-C₃ alkylsulfonyl,
- mono- or di-C₁-C₃ alkylaminosulfonyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
- C₁-C₇ alkyl,

- halogenated C₁-C₇ alkyl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (vi) heterocyclyl,
- or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkylthio substituted by carbocyclic aryl,
 - C₁-C₃ alkylthio substituted by halogenated carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - C₁-C₃ alkoxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkenylthio,
 - carbocyclic arylthio,
 - C₁-C₃ alkylsulfonyl,
 - carbocyclic arylsulfonyl,
 - halogenated carbocyclic arylsulfonyl,
 - carbocyclic arylsulfonyl substituted by C₁-C₄ alkyl,

- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
- halogen,
- nitro,
- C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
- C₁-C₃ alkyl,
- halogenated C₁-C₃ alkyl;

Y is -C(O)-;

wherein carbocyclic aryl is phenyl, naphthyl, anthranyl, or biphenyl;

carbocyclyl is 10,11-dihydro-5-oxo-dibenzo[a,d]cycloheptyl, 1-oxo-indanyl, 9H-fluorenyl, 9-oxo-fluorenyl, acenaphthyl, anthraquinonyl, C-fluoren-9-ylidene, indanyl, indenyl, 1,2,3,4-tetrahydro-naphthyl, or bicyclo[2.2.1]hepteny;

heterocyclyl is 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1,3-dioxo-isoindolyl, 1H-indolyl, 1H-pyrrolyl, 1-oxo-3H-isobenzofuranyl, 2,3-dihydro-benzo[1,4]dioxinyl, 2,3-dihydro-benzofuryl, 2,4-dihydro-3-oxo-pyrazolyl, 2H-benzopyranyl, 2-oxo-benzopyranyl, 2-oxo-pyrrolidinyl, 3,4-dihydro-2H-benzo[b][1,4]dioxepinyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, 4-oxo-3,4-dihydro-phthalazinyl, 4-oxo-benzopyranyl, 9,10,10-trioxo-thioxanthenyl, 9H-xanthenyl, azetidiny, benzimidazolyl, benzo[1,3]dioxolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, cinnolyl, furyl, imidazolyl, isoxazolyl, morpholino, morpholinyl, oxazolyl, oxolanyl, piperidyl, piridyl, pyrazolyl, pyridyl, pyrimidyl, pyrrolidyl, quinolyl, quinoxalyl, thiazolidyl, thiazolyl, thienyl, thiolanyl, tetrahydro-thienyl, benzofuranyl, or benzothiazolyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

3. A compound according to claim 2, wherein

R₁ represents

(i) C₁-C₁₀ alkyl,

C₁-C₁₀ alkyl substituted by substituent(s) independently selected from

- oxo,
- di-propylaminocarbonyl,
- methoxy substituted by carbocyclic aryl,
- methylcarbonyloxy,
- carbocyclic aryloxy,
- halogenated carbocyclic aryloxy,
- carbocyclic aryloxy substituted by nitro,
- heterocyclyloxy substituted by methyl,
- substituted heterocyclyl-ethylideneaminooxy,
- tert*-butoxycarbonylamino,
- carbocyclic arylcarbonylamino,
- C₁-C₂ alkylthio,
- C₁-C₂ alkylthio substituted by substituent(s) independently selected from
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by methoxy,
- carbocyclic arylthio,
- heterocyclylthio substituted by nitro,
- heterocyclylthio substituted by methyl,
- C₅-C₆ cycloalkyl,
- C₅-C₆ cycloalkenyl,
- carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - methyl,
 - methoxy,
 - ethenyl substituted by carbocyclic aryl substituted methylsulfinyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from

- oxo,
- carbocyclic aryl,
- heterocyclyl,
- C₁-C₄ alkoxy,
- halogenated C₁-C₄ alkoxy,
- C₁-C₄ alkoxy substituted by carbocyclic aryl,
- carbocyclic aryloxy,
- halogenated mono-carbocyclic arylaminocarbonyl,
- carbocyclic aryl,
- heterocyclyl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
- C₁-C₂ alkyl,
- C₁-C₂ substituted by carbocyclic aryl,
- methoxy,
- methoxy substituted by carbocyclic aryl,
- carbocyclic aryl,
- halogenated carbocyclic aryl,
- (ii) C₂-C₃ alkenyl substituted by substituent(s) independently selected from
- carbocyclic aryl,
- halogenated carbocyclic aryl,
- carbocyclic aryl substituted by nitro,
- (iii) C₃-C₆ cycloalkyl,
- C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from
- methyl substituted by oxo,
- methyl substituted by carbocyclic aryl,
- carbocyclic aryl,
- (iv) carbocyclyl,
- (v) carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
- halogen,
- hydroxy,

- cyano,
- nitro,
- C₁-C₉ alkyl,
- C₁-C₉ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryloxy,
 - C₁-C₇ alkoxy,
 - halogenated C₁-C₇ alkoxy,
 - C₁-C₇ alkoxy substituted by carbocyclic aryl,
 - methylcarbonyloxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methoxy,
 - amino,
 - di-methylamino,
 - propargynylcarbonylamino substituted by carbocyclic aryl,
 - carbocyclic arylsulfonylamino substituted by methyl,
 - (carbocyclic aryl)NHC(O)NH substituted by halogenated methoxy,
 - halogenated methylthio,
 - carbocyclic arylthio substituted by cyano,
 - di-propylamino sulfonyl,
 - mono- or di- ethylaminocarbonyl substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - heterocyclyl substituted by methyl,
 - heterocyclyl substituted by halogenated carbocyclic aryl,
- (vi) heterocyclyl,
or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,

- C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - methylthio substituted by halogenated carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - methoxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methyl,
 - C₁-C₃ alkylthio,
 - propenylthio,
 - carbocyclic arylthio,
 - C₁-C₃ alkylsulfonyl,
 - carbocyclic arylsulfonyl substituted by C₁-C₄ alkyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryl substituted by nitro,
 - heterocyclyl;

R₂ is methylamino or dimethylamino;

L is selected from Formula Va, VIIIa, or IXa;

wherein R₄ and R₅ are independently selected from H or C₁-C₃ alkyl;

wherein carbocyclic aryl is phenyl, naphthyl, anthranyl, or biphenyl;

carbocyclyl is 1-oxo-indanyl, 9-oxo-fluorenyl, indenyl, anthraquinonyl, C-fluoren-9-ylidene, 1,2,3,4-tetrahydro-naphthyl, or bicyclo[2.2.1]hepteny;

heterocyclyl is 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1,3-dioxo-isindolyl, 1*H*-indolyl, 1*H*-pyrrolyl, 1-oxo-3*H*-isobenzofuranyl, 2,3-dihydro-benzo[1,4]dioxinyl, 2,4-dihydro-3-oxo-pyrazolyl, 2*H*-benzopyranyl, 2-oxo-benzopyranyl, 3,4-dihydro-2*H*-benzo[b][1,4]dioxepinyl, 4-oxo-3,4-dihydro-phthalazinyl, 4-oxo-benzopyranyl, 9,10,10-trioxo-thioxanthenyl, 9*H*-xanthenyl, azetidiny, benzimidazolyl, benzo[1,3]dioxolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, furyl, imidazolyl, isoxazolyl, morpholino, morpholinyl, oxolanyl, piperidyl, piridyl, pyrazolyl, pyridyl, quinolyl,

quinoxalyl, thiazolidyl, thiazolyl, thienyl, thiolanyl, 2,3-dihydro-1-oxo-isindolyl, 2,3-dihydro-benzofuryl, 2-oxo-pyrrolidinyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, cinnolyl, pyrimidyl, pyrrolidyl, tetrahydro-thienyl, benzofuranyl, or benzothiazolyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

4. A compound according to claim 3, wherein

R₁ represents

(i) C₁-C₁₀ alkyl substituted by substituent(s) independently selected from

- oxo,
- di-propylaminocarbonyl,
- methoxy substituted by carbocyclic aryl,
- methylcarbonyloxy,
- carbocyclic aryloxy,
- halogenated carbocyclic aryloxy,
- carbocyclic aryloxy substituted by nitro,
- heterocyclyloxy substituted by methyl,
- substituted heterocyclyl-ethylideneaminooxy,
- tert*-butoxycarbonylamino,
- carbocyclic arylcarbonylamino,
- C₁-C₂ alkylthio,
- C₁-C₂ alkylthio substituted by substituent(s) independently selected from
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by methoxy,
 - carbocyclic arylthio,
 - heterocyclylthio substituted by nitro,
 - heterocyclylthio substituted by methyl,
- C₅-C₆ cycloalkenyl,
- carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - methyl,
 - methoxy,

- ethenyl substituted by carbocyclic aryl substituted methylsulfinyl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - oxo,
 - carbocyclic aryl,
 - heterocyclyl,
 - C₁-C₄ alkoxy,
 - halogenated C₁-C₄ alkoxy,
 - C₁-C₄ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryloxy,
 - halogenated mono-carbocyclic arylaminocarbonyl,
 - carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₂ alkyl,
 - C₁-C₂ substituted by carbocyclic aryl,
 - methoxy,
 - methoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (ii) C₂-C₃ alkenyl substituted by substituent(s) independently selected from
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by nitro,
- (iii) C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from
 - methyl substituted by oxo,
 - methyl substituted by carbocyclic aryl,
 - carbocyclic aryl,

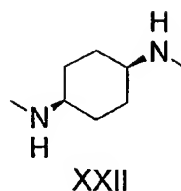
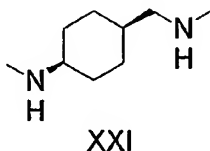
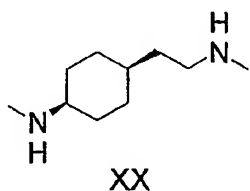
(iv) carbocyclyl,

(v) carbocyclic aryl substituted by substituent(s) independently selected from

- halogen,
- hydroxy,
- cyano,
- nitro,
- C₁-C₉ alkyl,
- C₁-C₉ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryloxy,
 - C₁-C₇ alkoxy,
 - halogenated C₁-C₇ alkoxy,
 - C₁-C₇ alkoxy substituted by carbocyclic aryl,
 - methylcarbonyloxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methoxy,
 - amino,
 - di-methylamino,
 - propargynylcarbonylamino substituted by carbocyclic aryl,
 - carbocyclic arylsulfonylamino substituted by methyl,
 - (carbocyclic aryl)NHC(O)NH substituted by halogenated methoxy,
 - halogenated methylthio,
 - carbocyclic arylthio substituted by cyano,
 - di-propylamino sulfonyl,
 - mono- or di- ethylaminocarbonyl substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - heterocyclyl substituted by methyl,
 - heterocyclyl substituted by halogenated carbocyclic aryl,
- (vi) or heterocyclyl substituted by substituent(s) independently selected from

- halogen,
- nitro,
- C₁-C₄ alkyl,
- C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - methylthio substituted by halogenated carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - methoxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methyl,
 - C₁-C₃ alkylthio,
 - propenylthio,
 - carbocyclic arylthio,
 - C₁-C₃ alkylsulfonyl,
 - carbocyclic arylsulfonyl,
 - carbocyclic arylsulfonyl substituted by C₁-C₄ alkyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryl substituted by nitro,
 - heterocyclyl;

L is selected from Formula XX - XXII;



wherein carbocyclic aryl is phenyl, naphthyl, or biphenyl;

carbocyclyl is 1-oxo-indanyl, 9-oxo-fluorenyl, indenyl, anthraquinonyl, C-fluoren-

9-ylidene, 1,2,3,4-tetrahydro-naphthyl, or bicyclo[2.2.1]hepteny;

heterocyclyl is 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1*H*-indolyl, 1*H*-pyrrolyl, 2,4-dihydro-3-oxo-pyrazolyl, 2*H*-benzopyranyl, 4-oxo-benzopyranyl, azetidiny, benzo[b]thienyl, furyl, isoxazolyl, morpholinyl, piperidyl, piridyl, pyrazolyl, pyridyl, quinolyl, thiazolidyl, thiazolyl, thienyl, thiolanyl, 2,3-dihydro-1-oxo-isoindolyl, 2,3-dihydro-benzofuryl, 2-oxo-benzopyranyl, 2-oxo-pyrrolidinyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, 9*H*-xanthenyl, cinnolyl, imidazolyl, morpholino, pyrimidyl, pyrrolidyl, tetrahydro-thienyl, benzofuranyl, or benzothiazolyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

5. A compound according to claim 4, wherein

R₁ represents

(i) C₁-C₅ alkyl substituted by substituent(s) independently selected from

- oxo,
- di-propylaminocarbonyl,
- methoxy substituted by carbocyclic aryl,
- methylcarbonyloxy,
- carbocyclic aryloxy,
- halogenated carbocyclic aryloxy,
- carbocyclic aryloxy substituted by nitro,
- heterocyclyloxy substituted by methyl,
- substituted heterocyclyl-ethylideneaminooxy,
- tert*-butoxycarbonylamino,
- carbocyclic arylcarbonylamino,
- C₁-C₂ alkylthio,
- C₁-C₂ alkylthio substituted by substituent(s) independently selected from
- halogenated carbocyclic aryl,
- carbocyclic aryl substituted by methoxy,
- carbocyclic arylthio,
- hetrocyclylthio substituted by nitro,
- hetrocyclylthio substituted by methyl,

- cyclohexenyl,
- carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - methyl,
 - methoxy,
 - ethenyl substituted by carbocyclic aryl substituted methylsulfinyl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - oxo,
 - carbocyclic aryl,
 - heterocyclyl,
 - C₁-C₂ alkoxy,
 - halogenated C₁-C₂ alkoxy,
 - C₁-C₂ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryloxy,
 - halogenated mono-carbocyclic arylaminocarbonyl,
 - carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₂ alkyl,
 - C₁-C₂ substituted by carbocyclic aryl,
 - methoxy,
 - methoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (ii) C₂-C₃ alkenyl substituted by substituent(s) independently selected from
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,

- carbocyclic aryl substituted by nitro,
- (iii) C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from
 - methyl substituted by oxo,
 - methyl substituted by carbocyclic aryl,
 - carbocyclic aryl,
- (iv) carbocyclyl,
- (v) carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₂ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - oxo,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryloxy,
 - C₁-C₂ alkoxy,
 - halogenated C₁-C₂ alkoxy,
 - C₁-C₂ alkoxy substituted by carbocyclic aryl,
 - methylcarbonyloxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methoxy,
 - amino,
 - di-methylamino,
 - propargynylcarbonylamino substituted by carbocyclic aryl,
 - carbocyclic arylsulfonylamino substituted by methyl,
 - (carbocyclic aryl)NHC(O)NH substituted by halogenated methoxy,
 - halogenated methylthio,
 - carbocyclic arylthio substituted by cyano,
 - di-propylamino sulfonyl,

- mono- or di- ethylaminocarbonyl substituted by carbocyclic aryl,
- carbocyclic aryl,
- heterocyclyl substituted by methyl,
- heterocyclyl substituted by halogenated carbocyclic aryl,
- (vi) or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - methylthio substituted by halogenated carbocyclic aryl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - methoxy,
 - carbocyclic aryloxy,
 - carbocyclic aryloxy substituted by methyl,
 - C₁-C₃ alkylthio,
 - propenylthio,
 - carbocyclic arylthio,
 - C₁-C₃ alkylsulfonyl,
 - carbocyclic arylsulfonyl,
 - carbocyclic arylsulfonyl substituted by methyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - carbocyclic aryl substituted by methyl,
 - carbocyclic aryl substituted by nitro,
 - heterocyclyl;

wherein carbocyclic aryl is phenyl, naphthyl, or biphenyl;

carbocyclyl is 1-oxo-indanyl, indenyl, 9-oxo-fluorenyl, 1,2,3,4-tetrahydro-naphthyl, or bicyclo[2.2.1]hepteny;

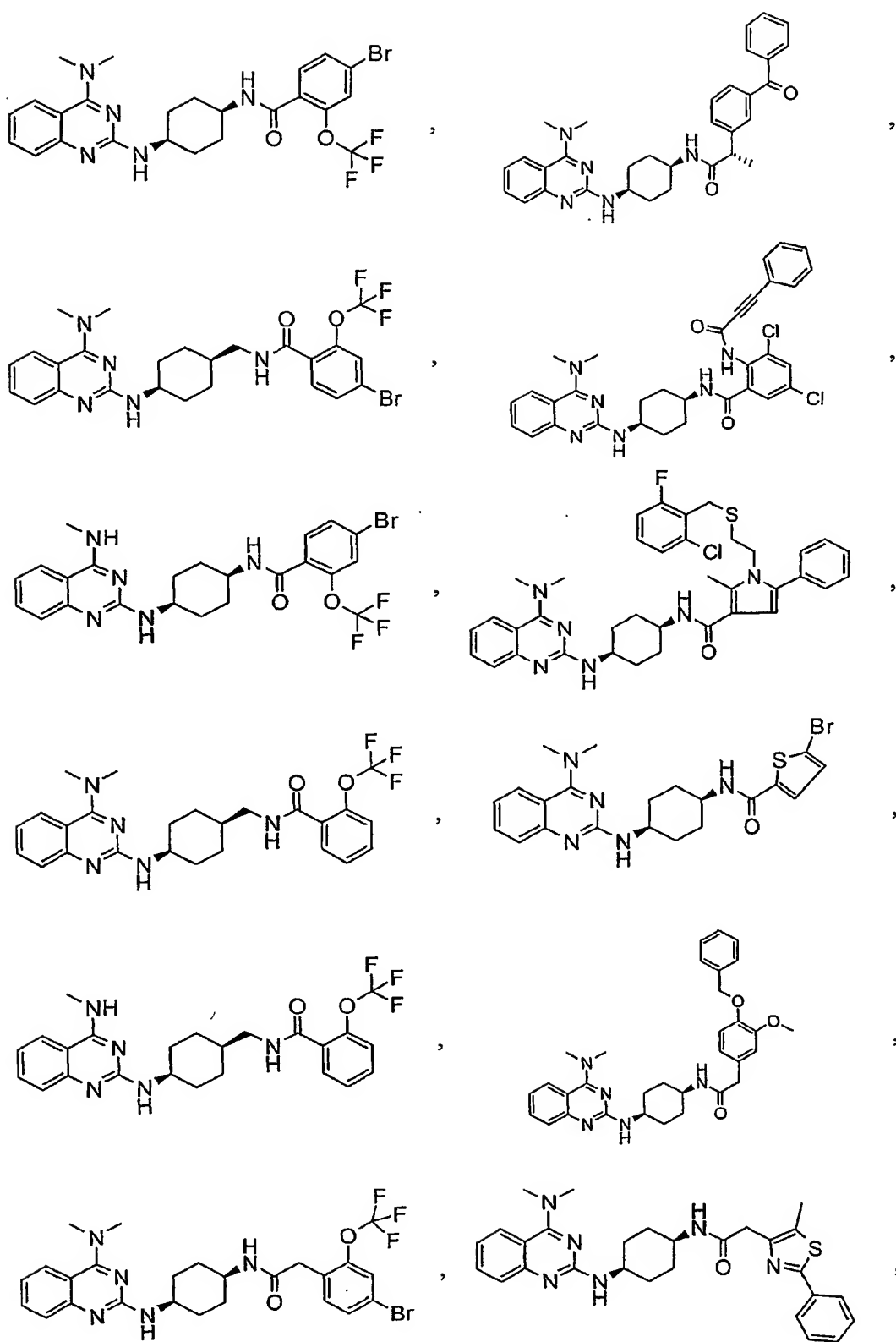
heterocyclyl is 1*H*-indolyl, 2,4-dihydro-3-oxo-pyrazolyl, furyl, pyrazolyl, pyridyl,

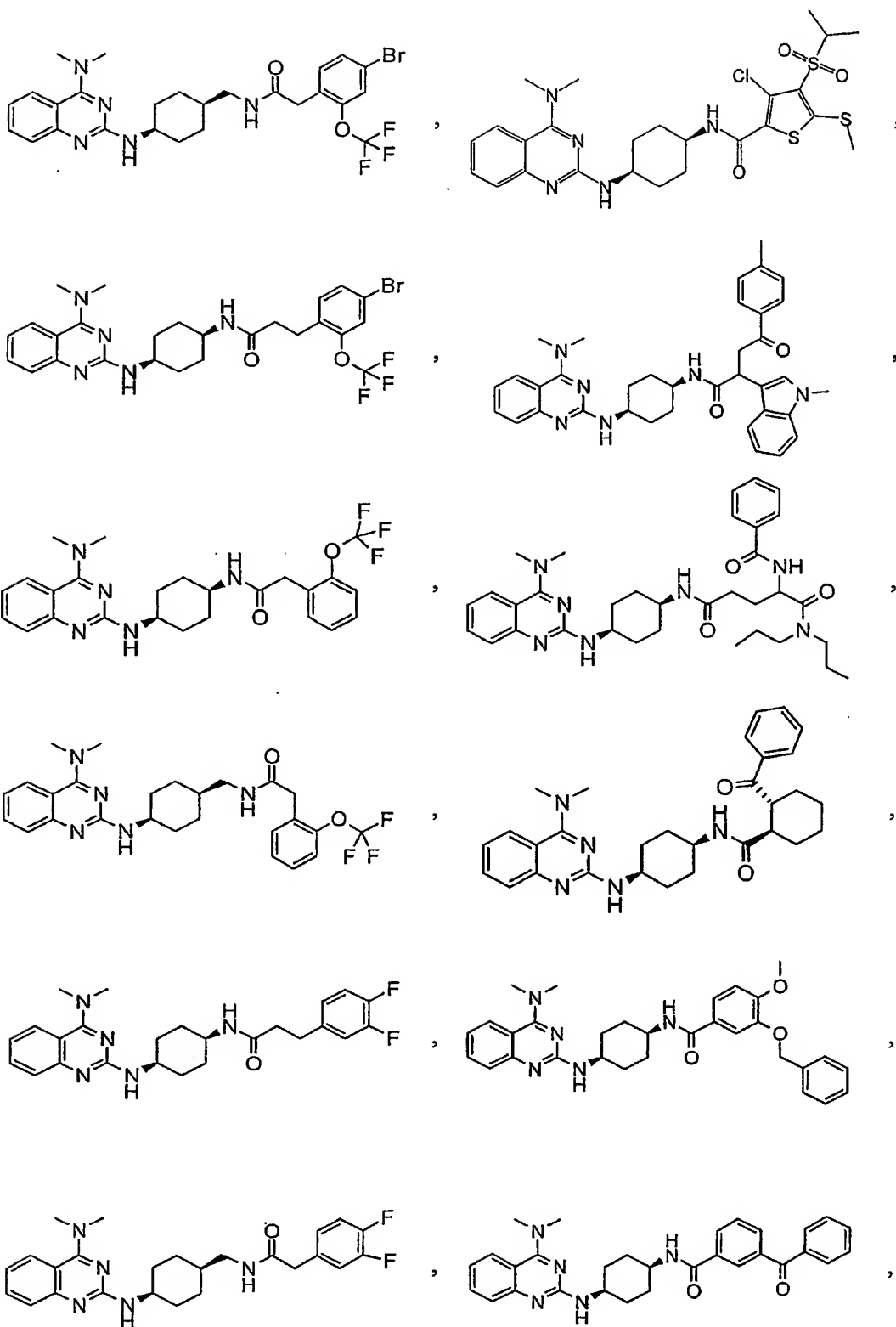
thienyl, 1,2,3-triazolyl, 1*H*-pyrrolyl, 2,3-dihydro-1-oxo-isoindolyl, 2,3-dihydro-benzofuryl, 2*H*-benzopyranyl, 2-oxo-benzopyranyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, imidazolyl, isoxazolyl, morpholino, morpholinyl, pyrazolyl, pyrimidyl, quinolyl, thiazolyl, tetrahydro-thienyl, benzofuranyl, or benzothiazolyl;

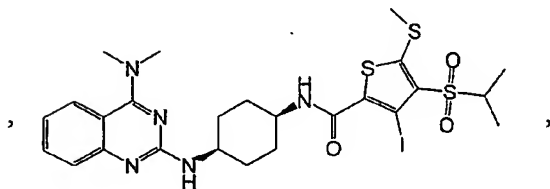
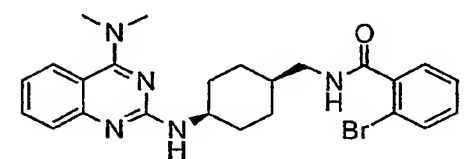
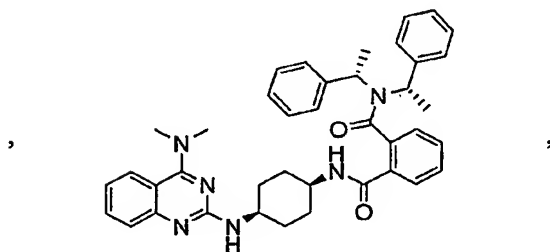
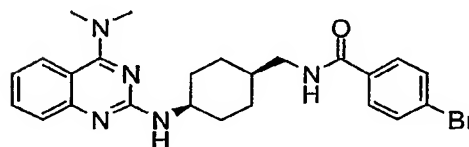
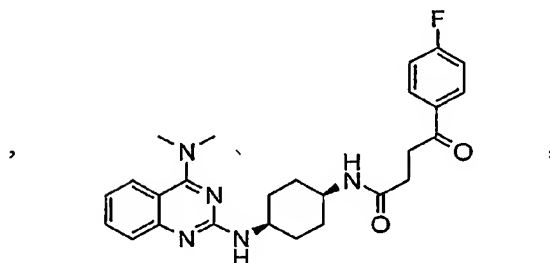
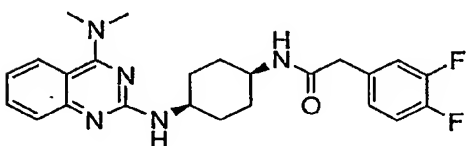
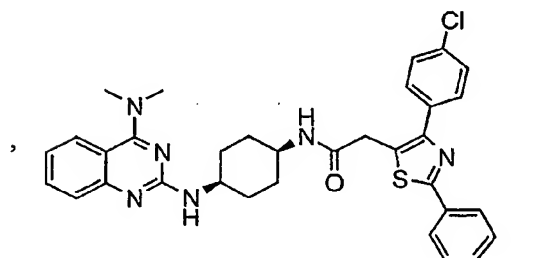
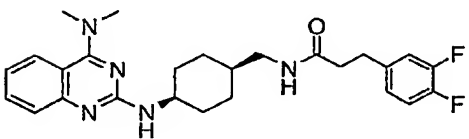
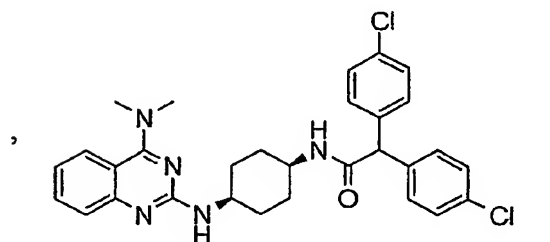
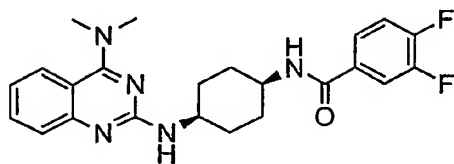
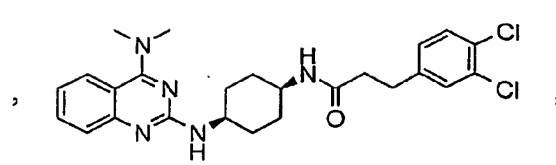
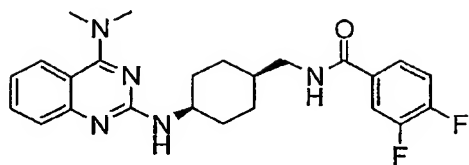
halogen is fluoro, chloro, bromo, or iodo;

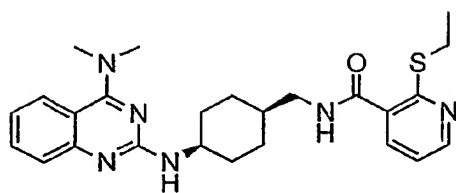
or a salt thereof.

6. A compound according to claim 5 of Formula I selected from the group consisting of

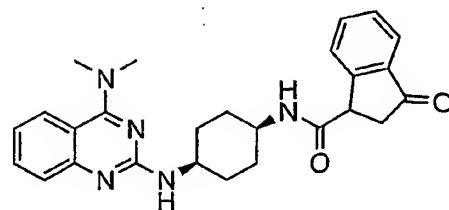




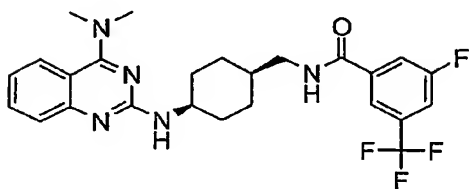




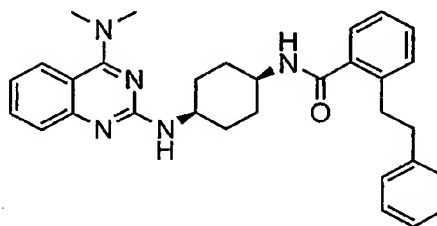
,



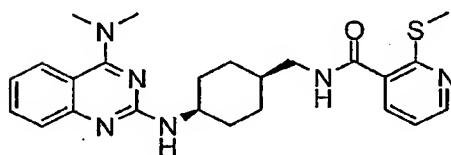
,



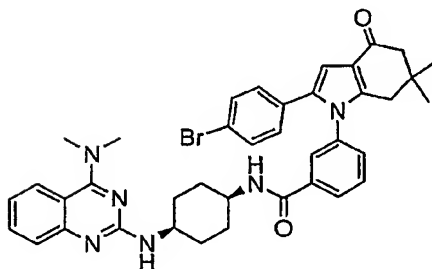
,



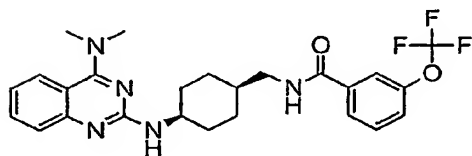
,



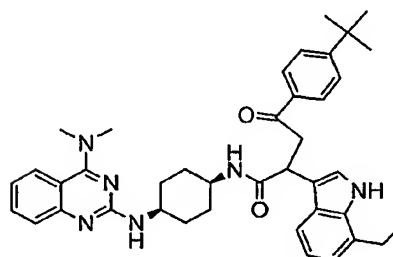
,



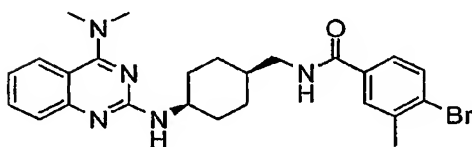
,



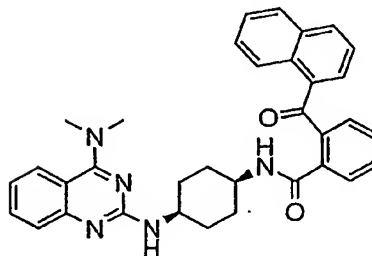
,



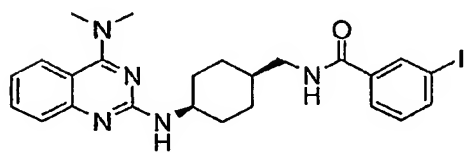
,



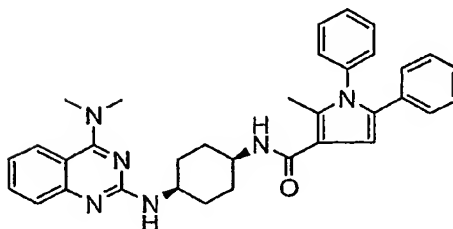
,



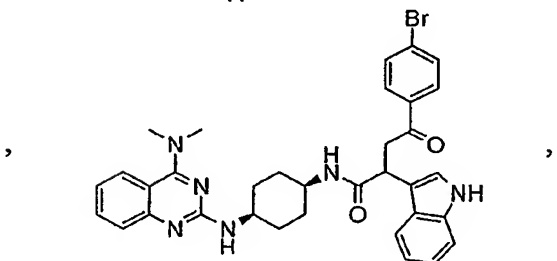
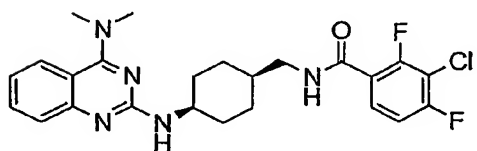
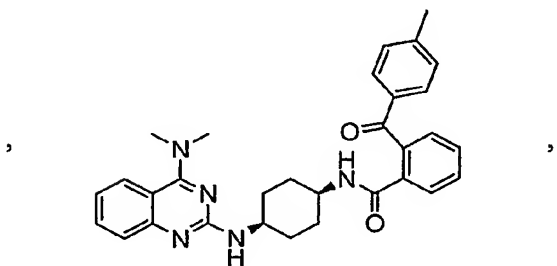
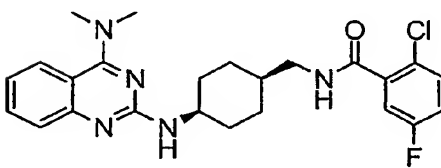
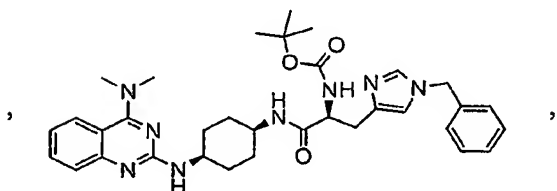
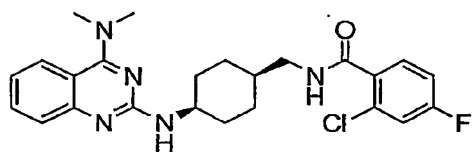
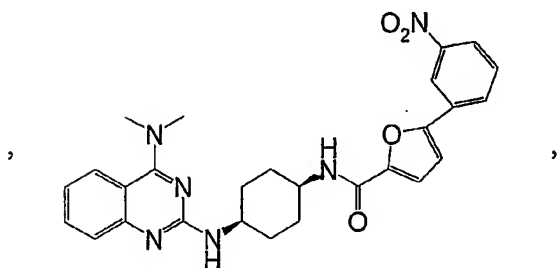
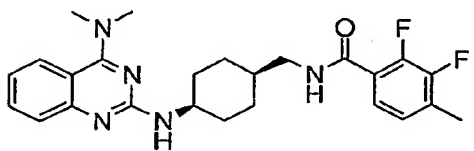
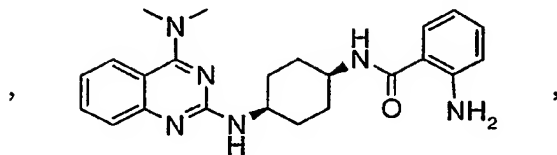
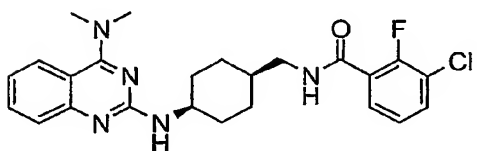
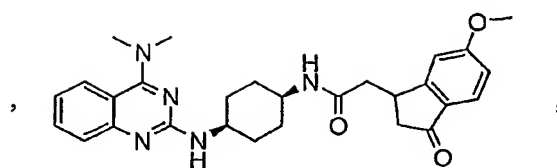
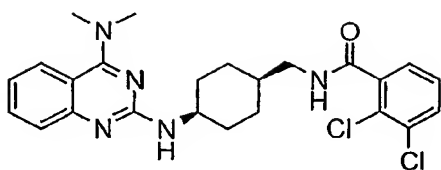
,

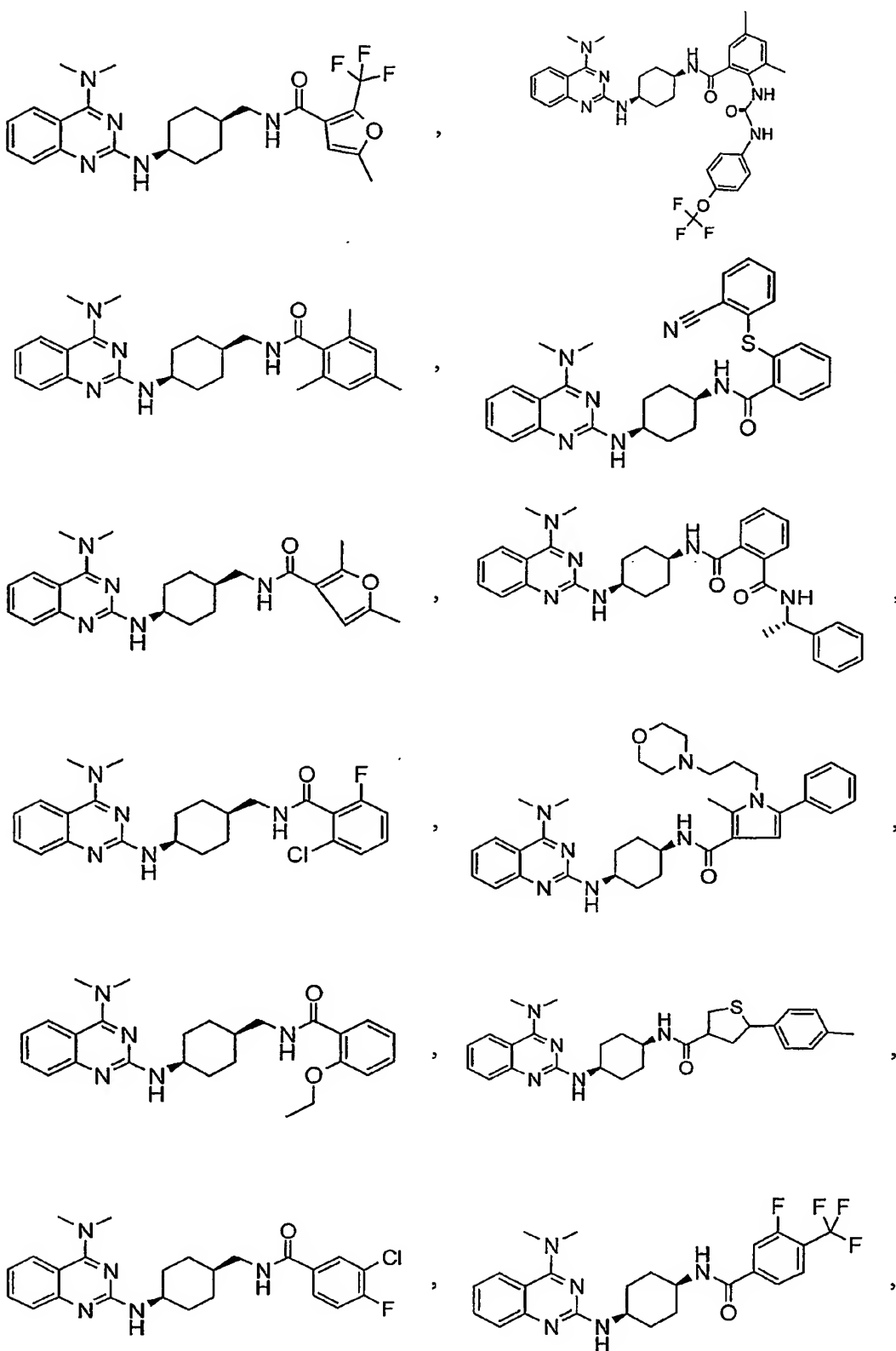


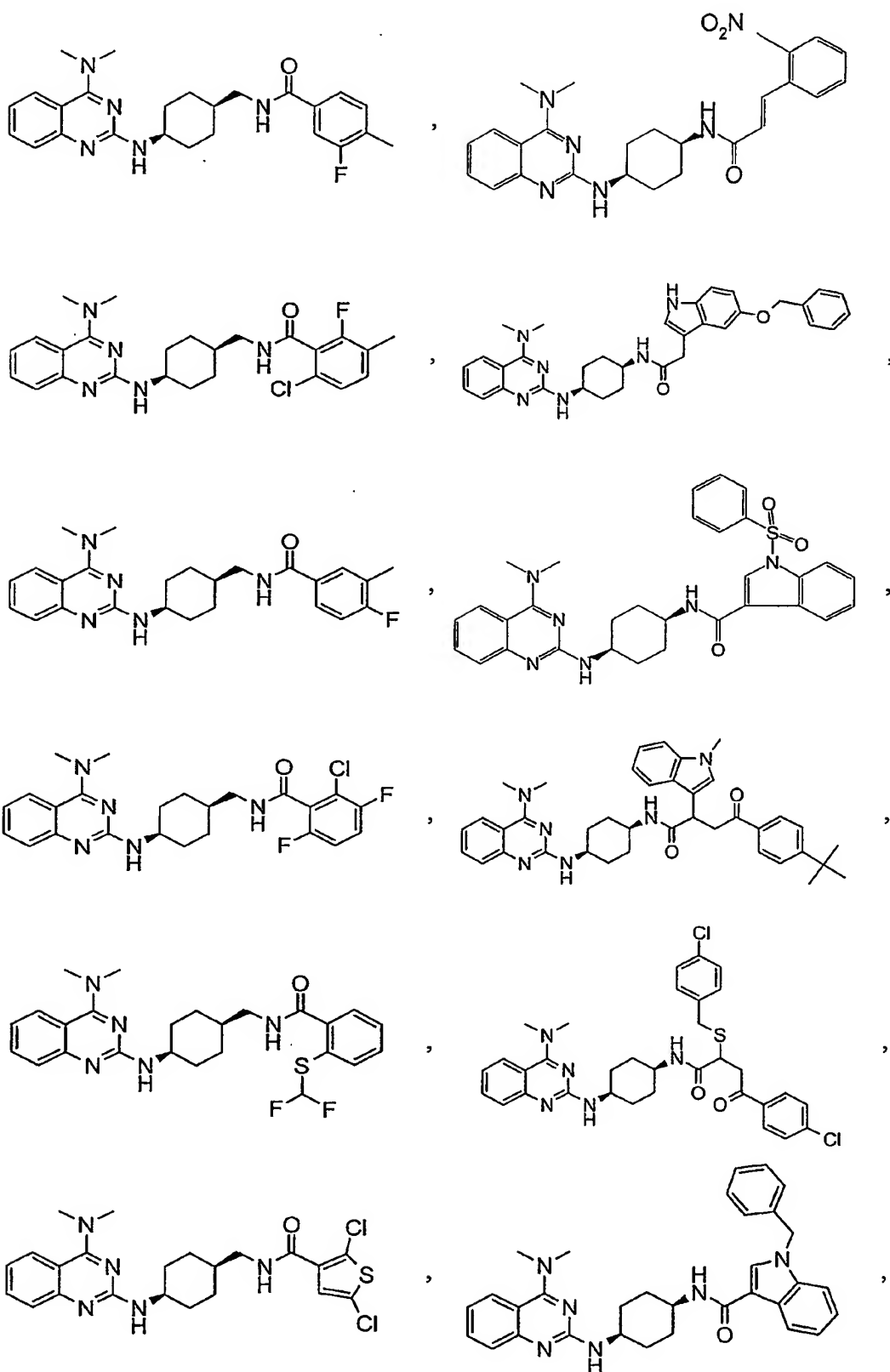
,

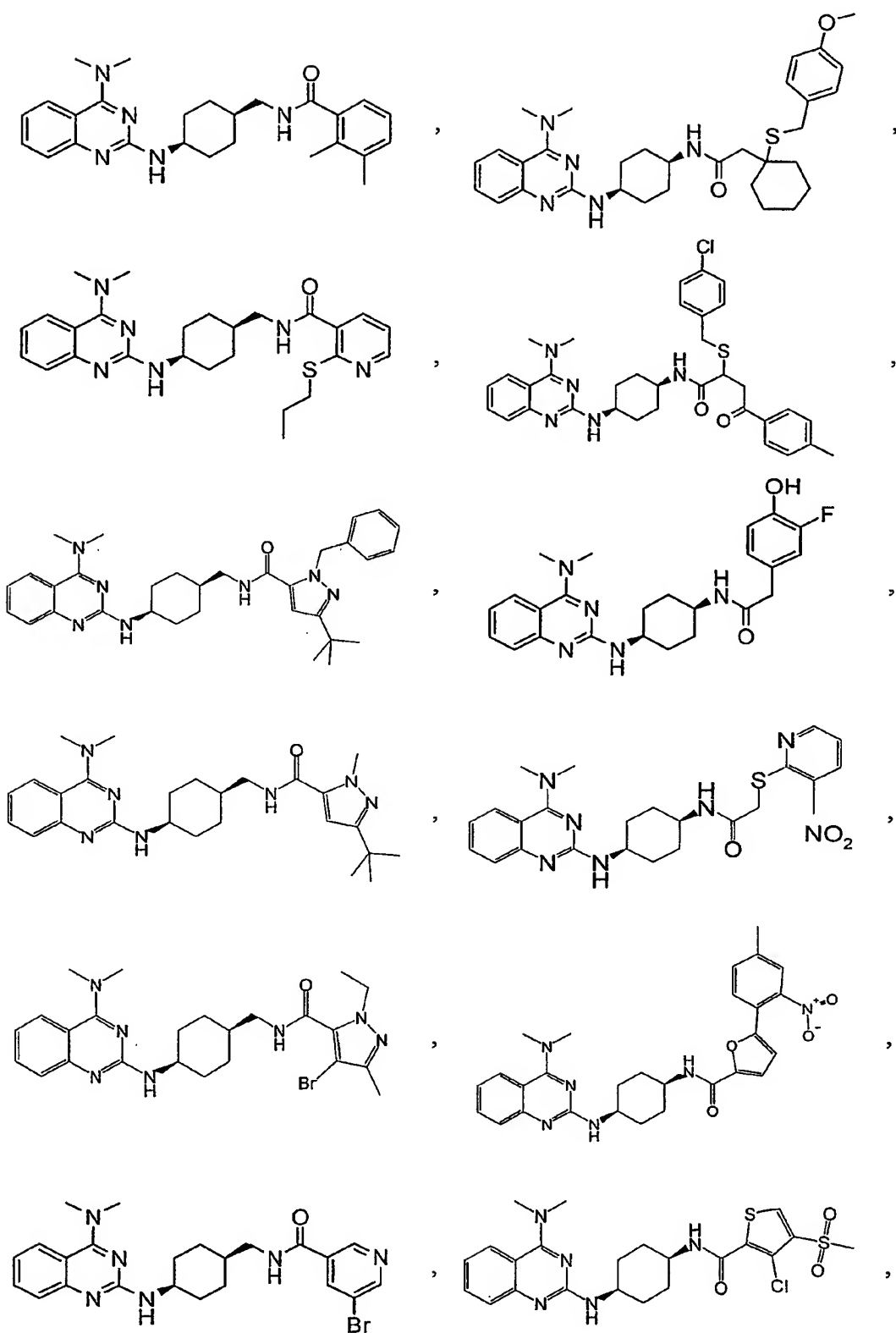


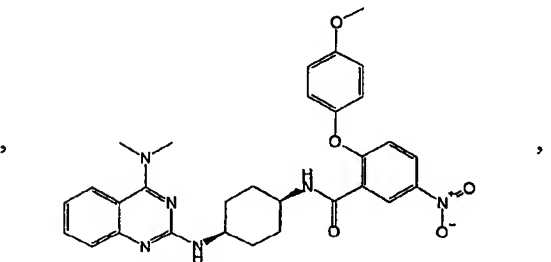
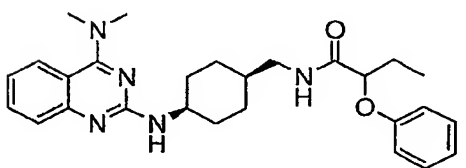
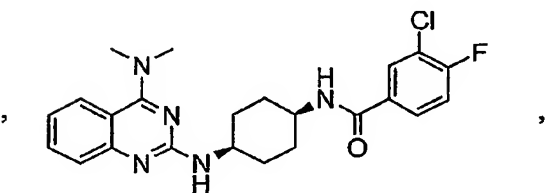
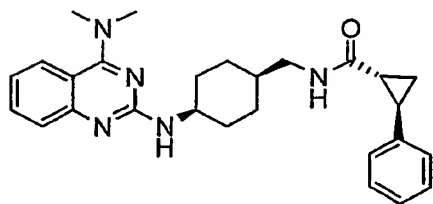
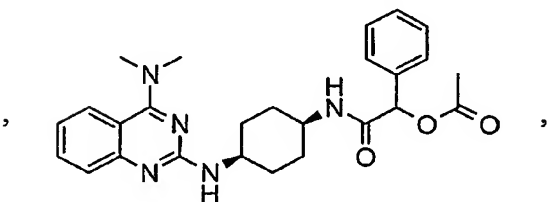
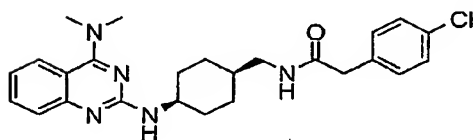
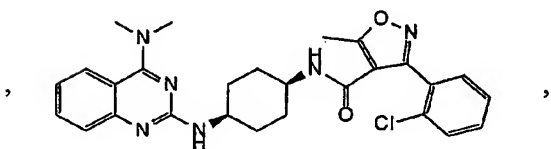
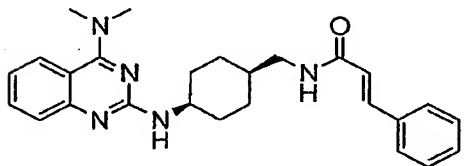
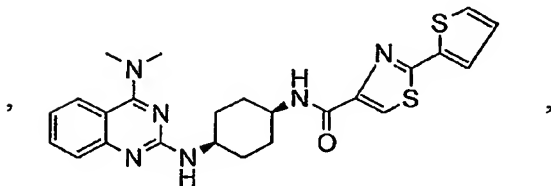
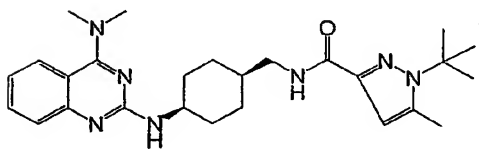
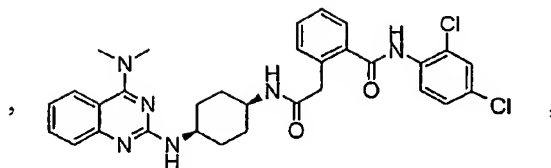
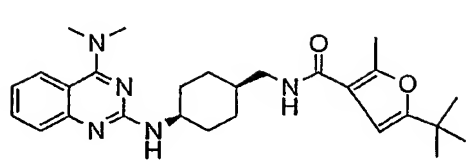
,

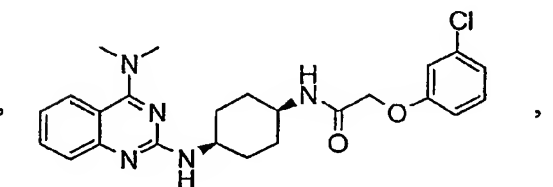
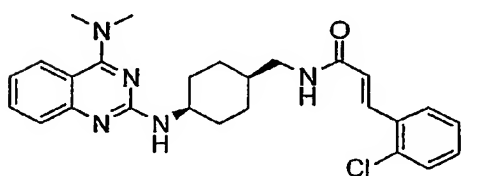
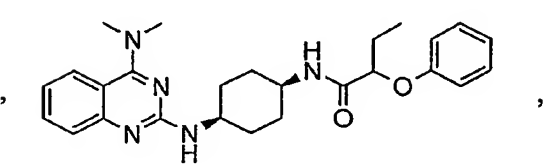
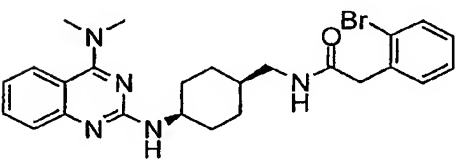
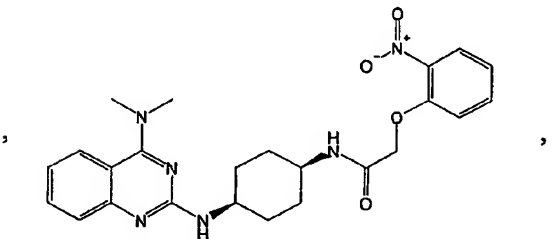
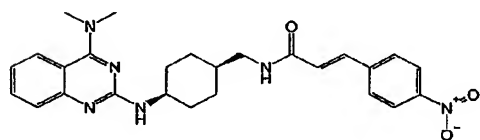
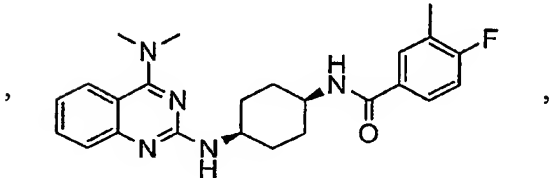
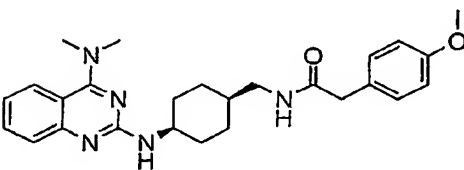
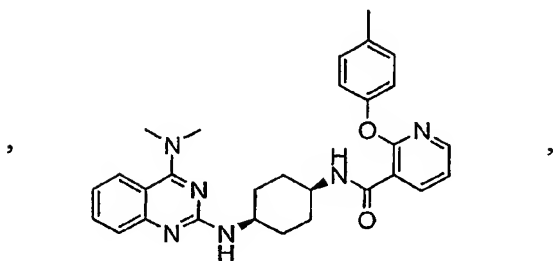
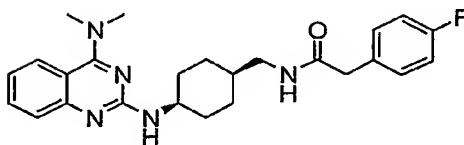
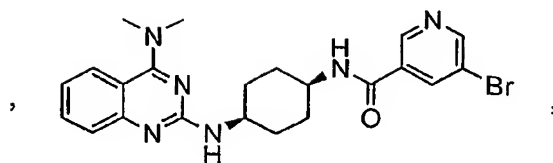
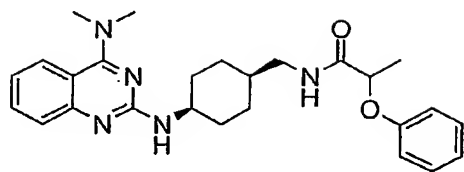


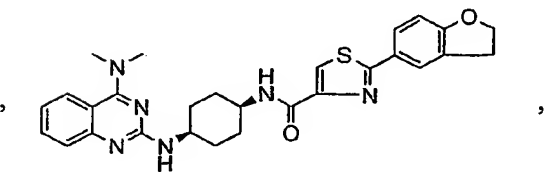
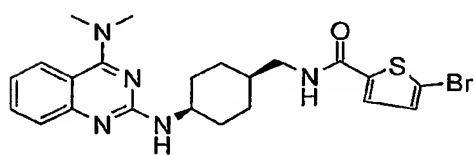
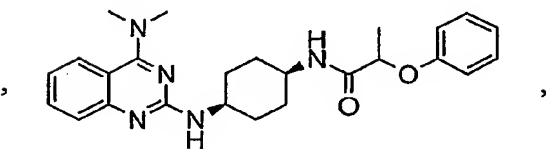
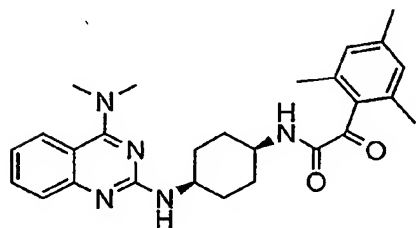
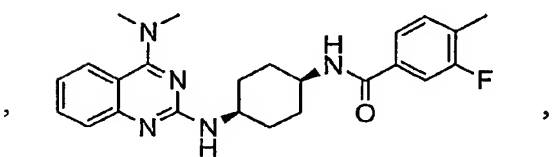
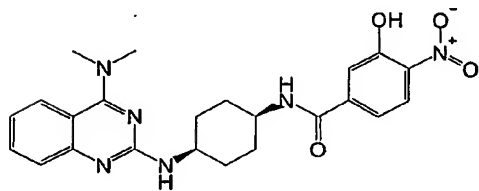
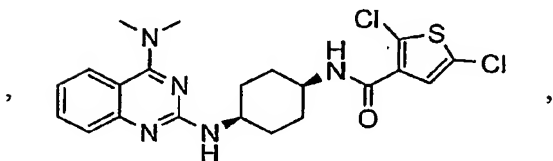
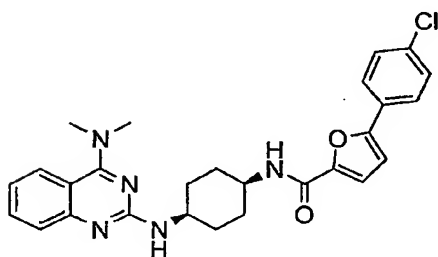
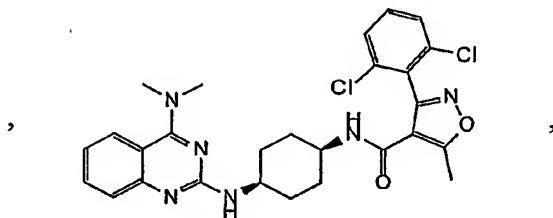
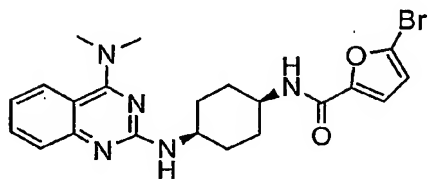
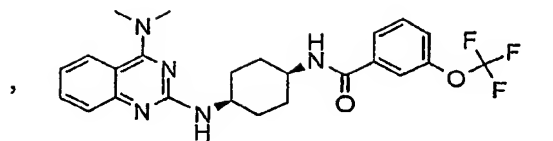
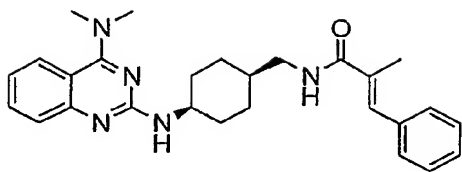


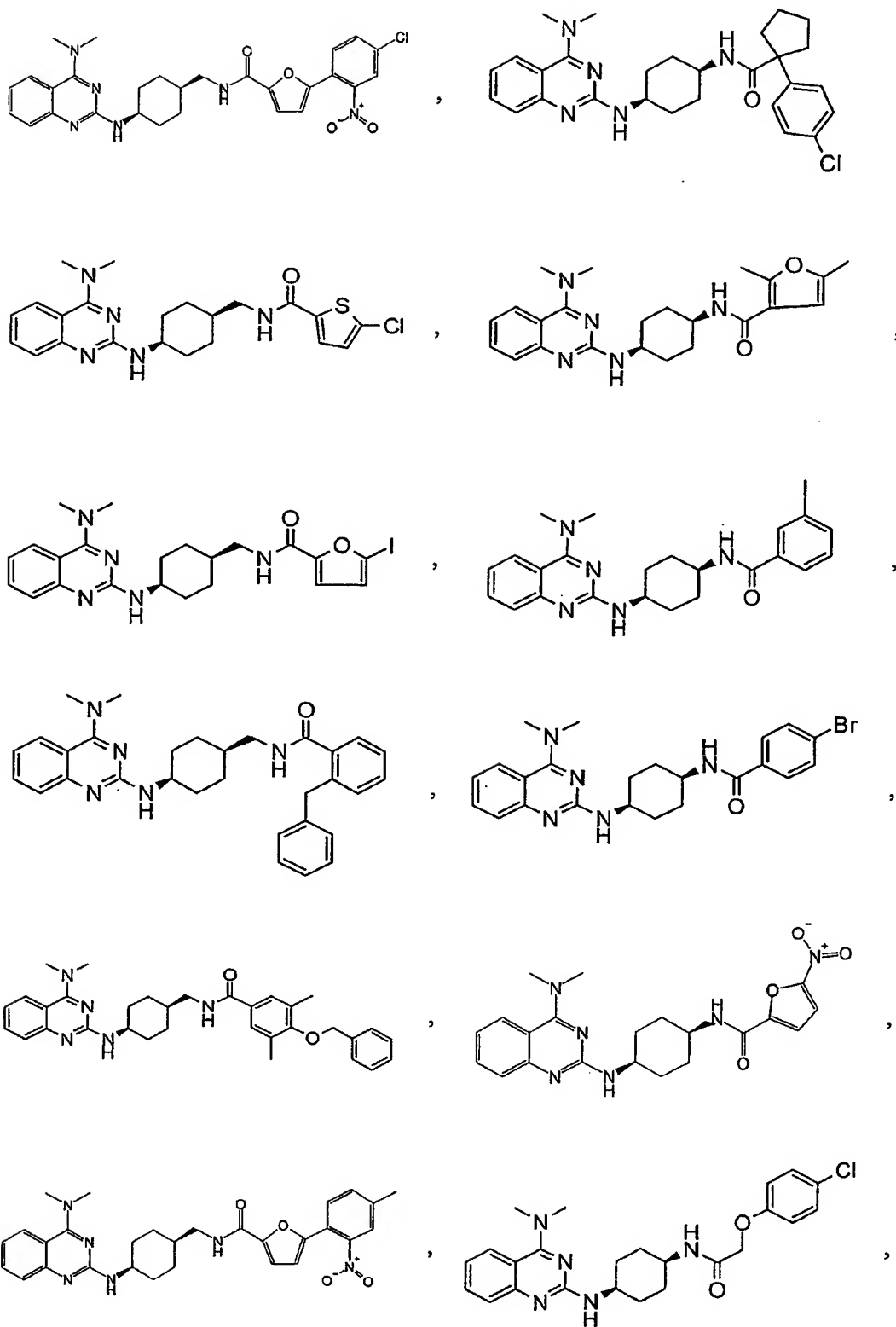


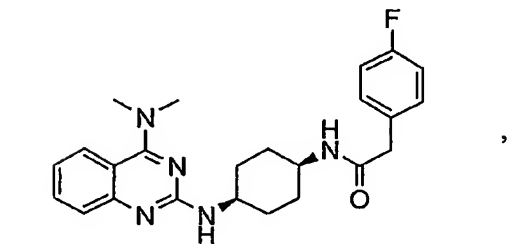
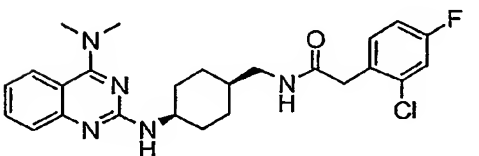
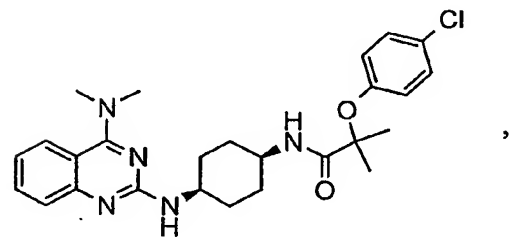
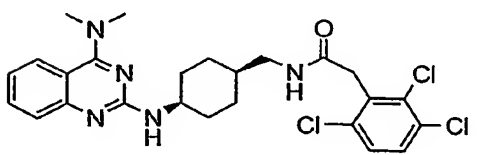
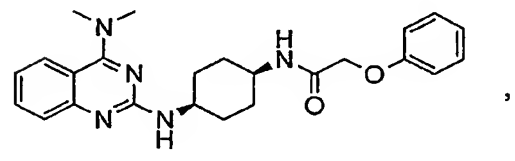
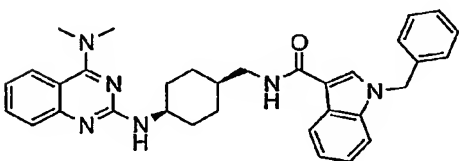
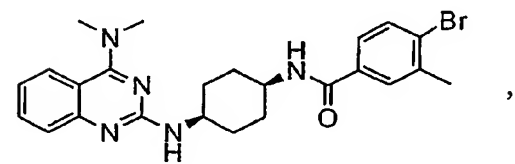
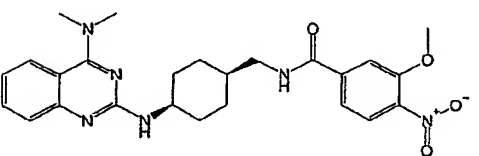
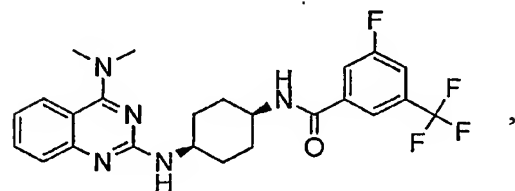
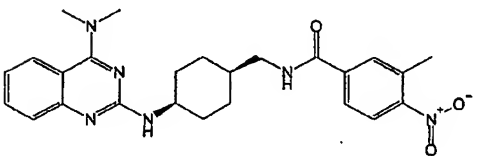
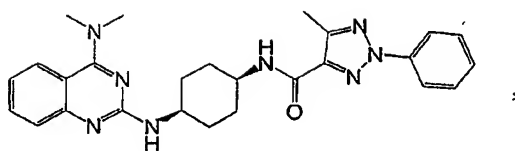
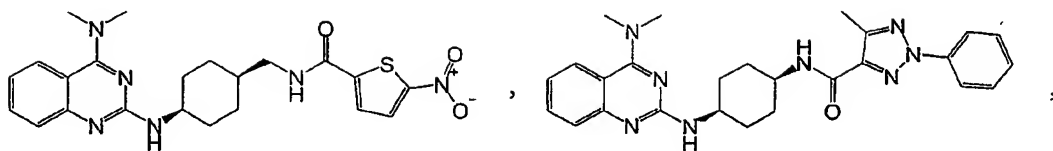


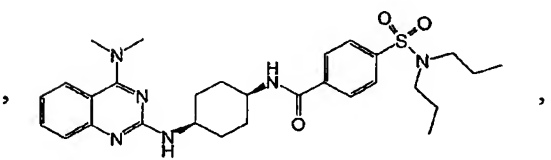
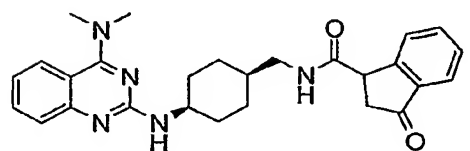
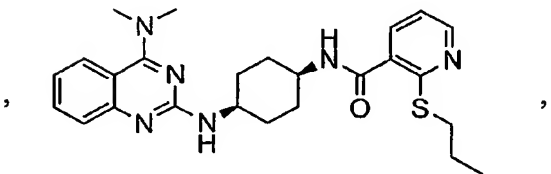
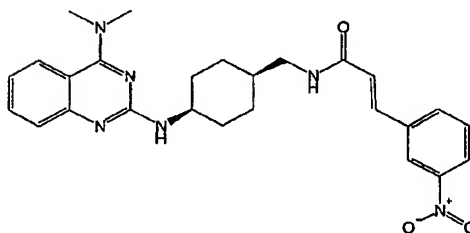
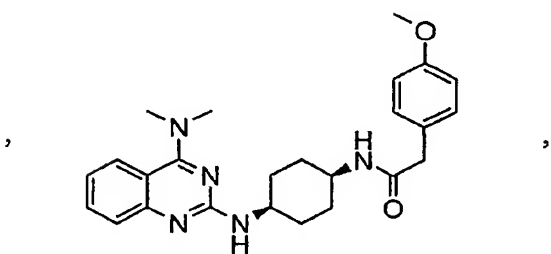
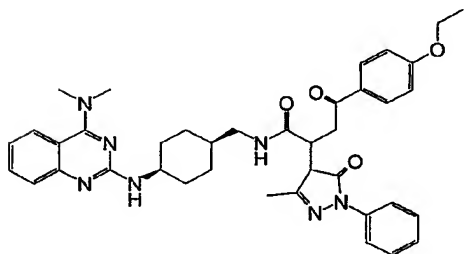
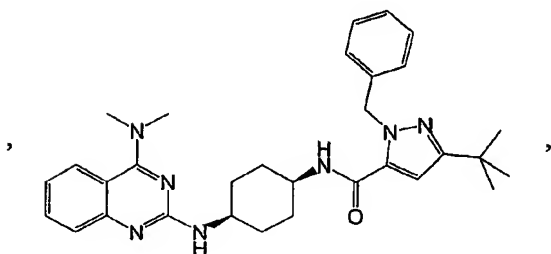
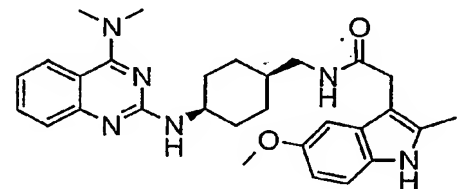
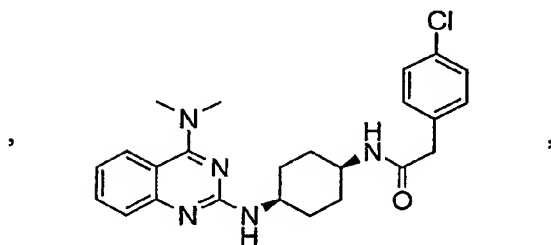
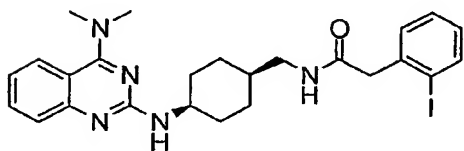
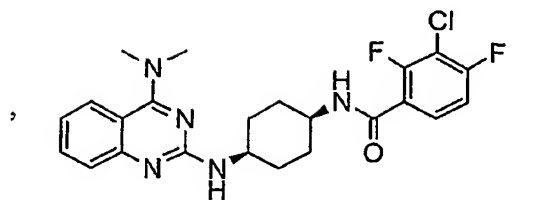
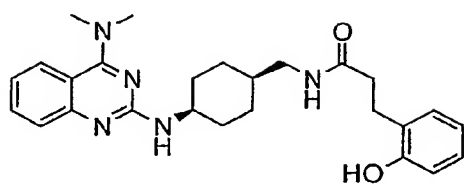


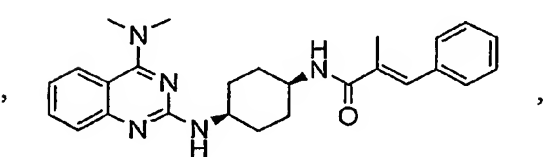
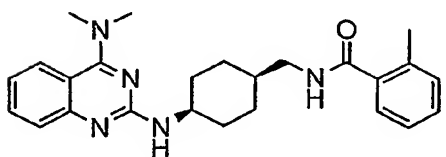
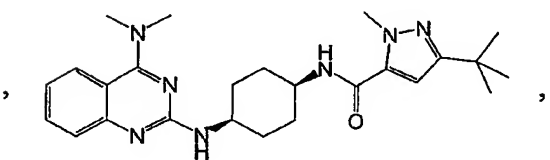
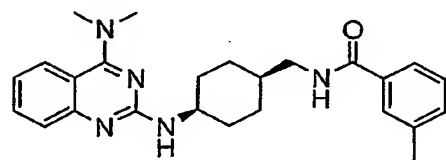
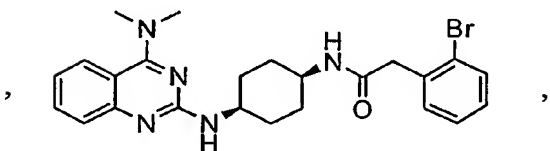
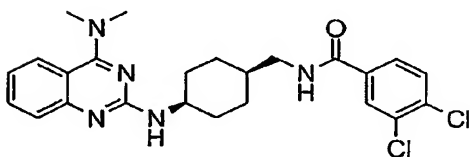
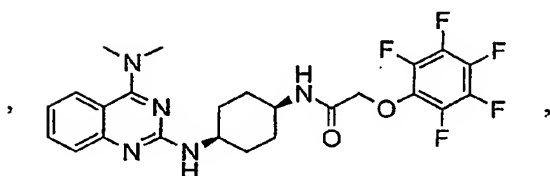
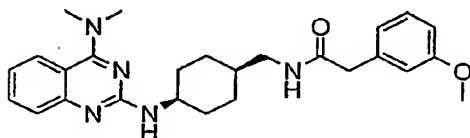
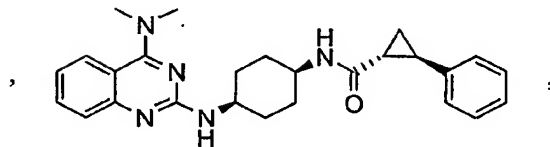
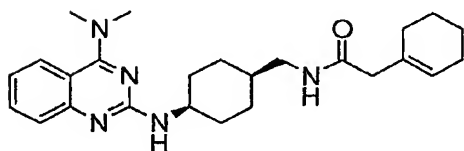
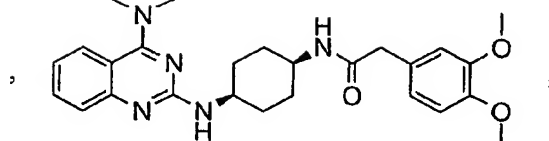
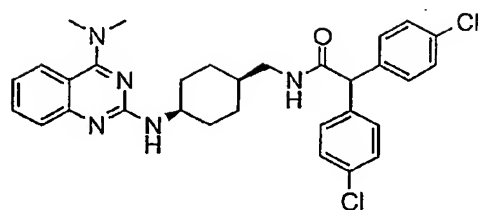


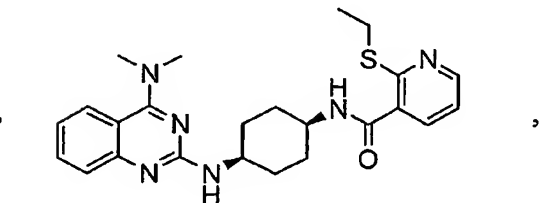
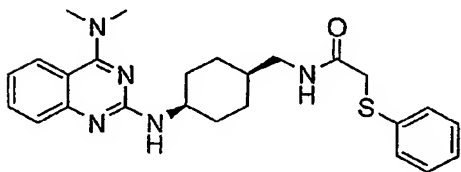
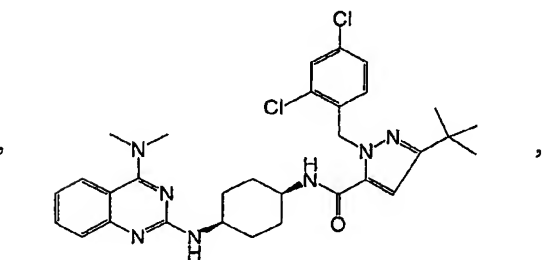
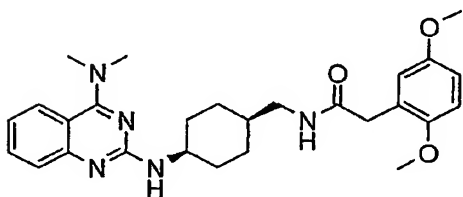
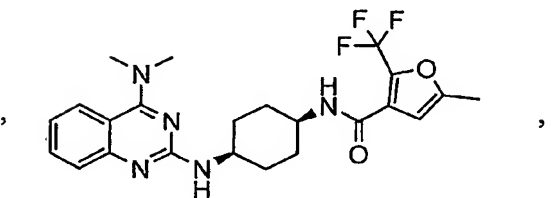
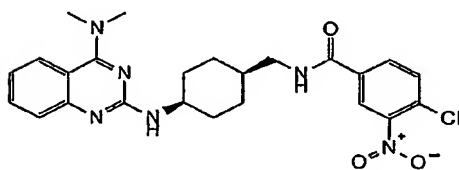
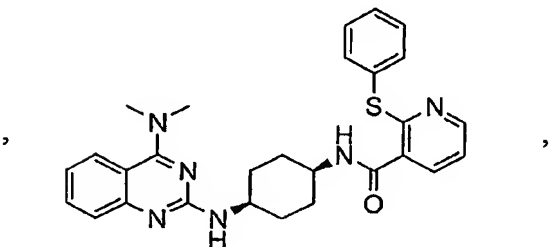
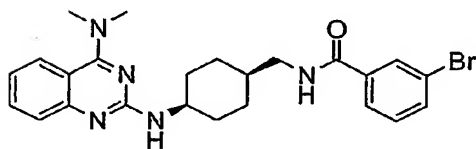
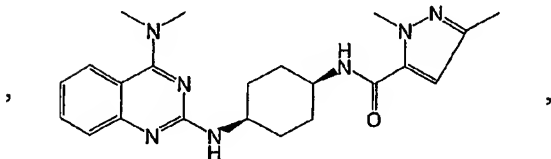
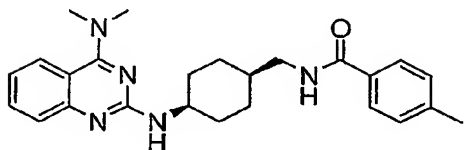
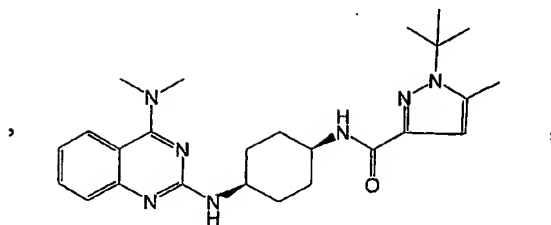
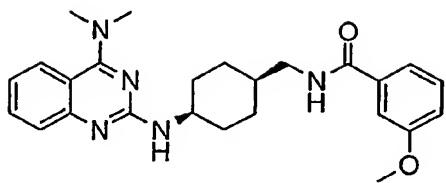


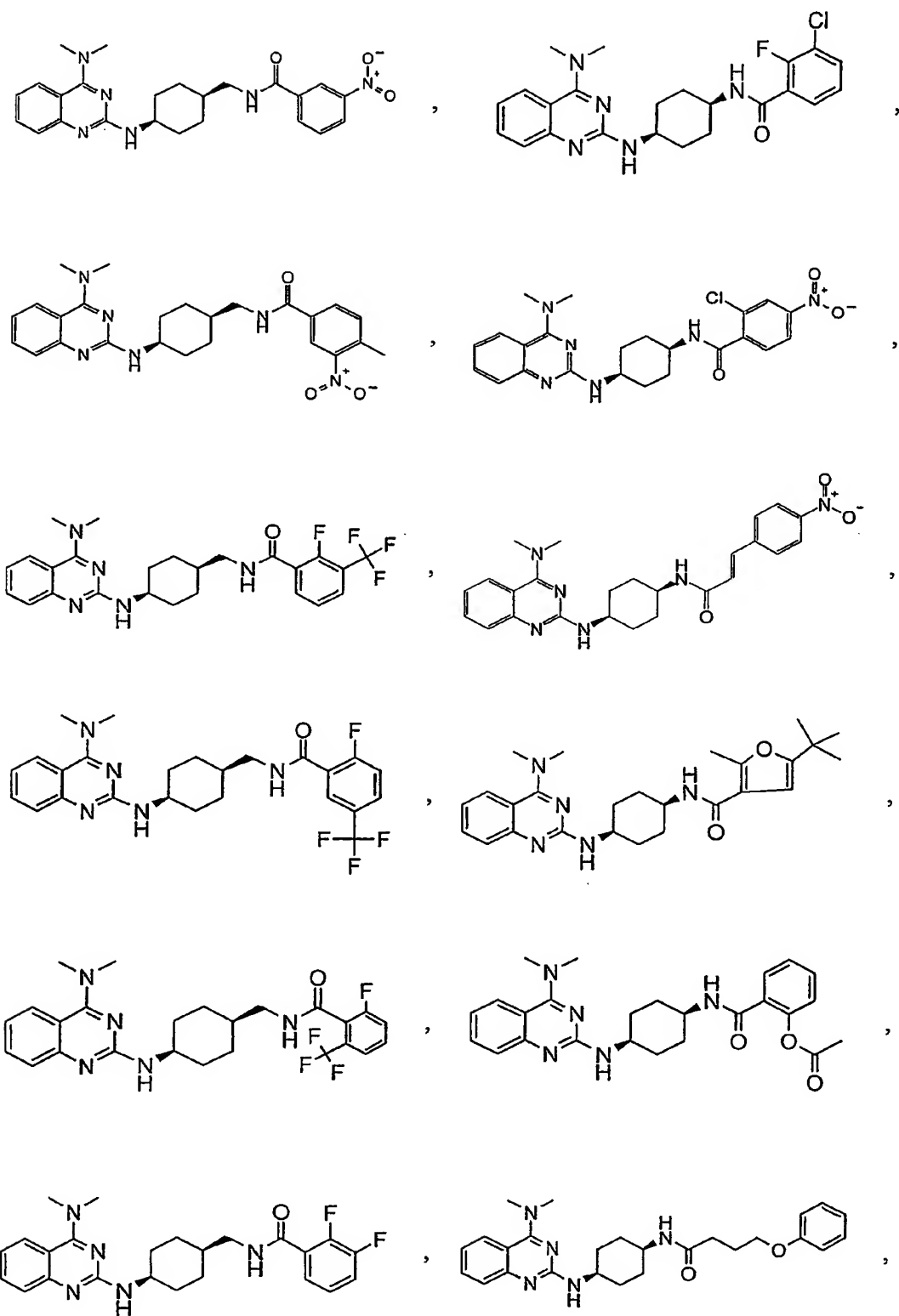


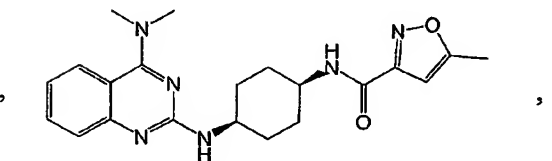
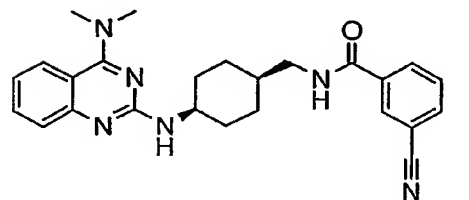
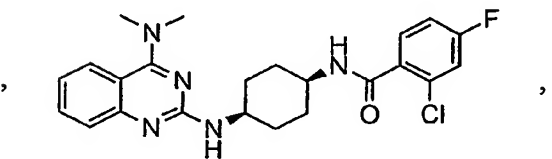
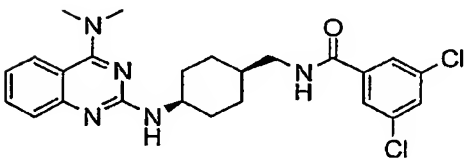
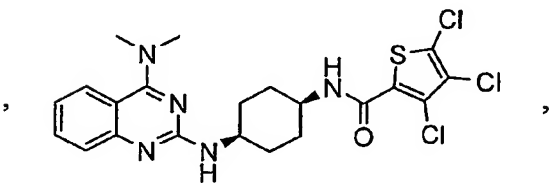
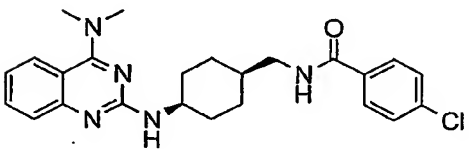
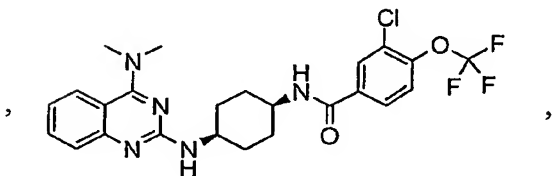
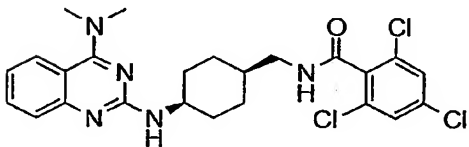
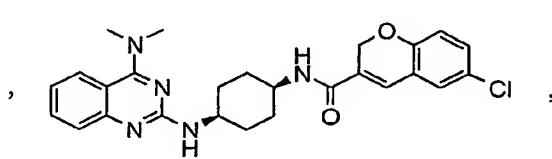
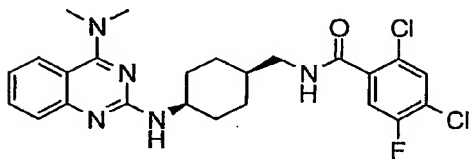
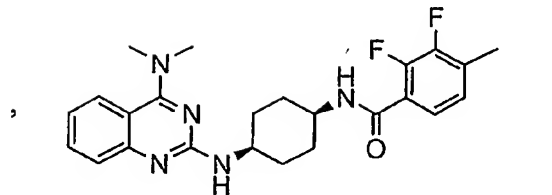
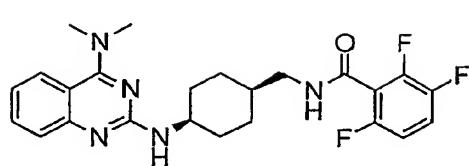


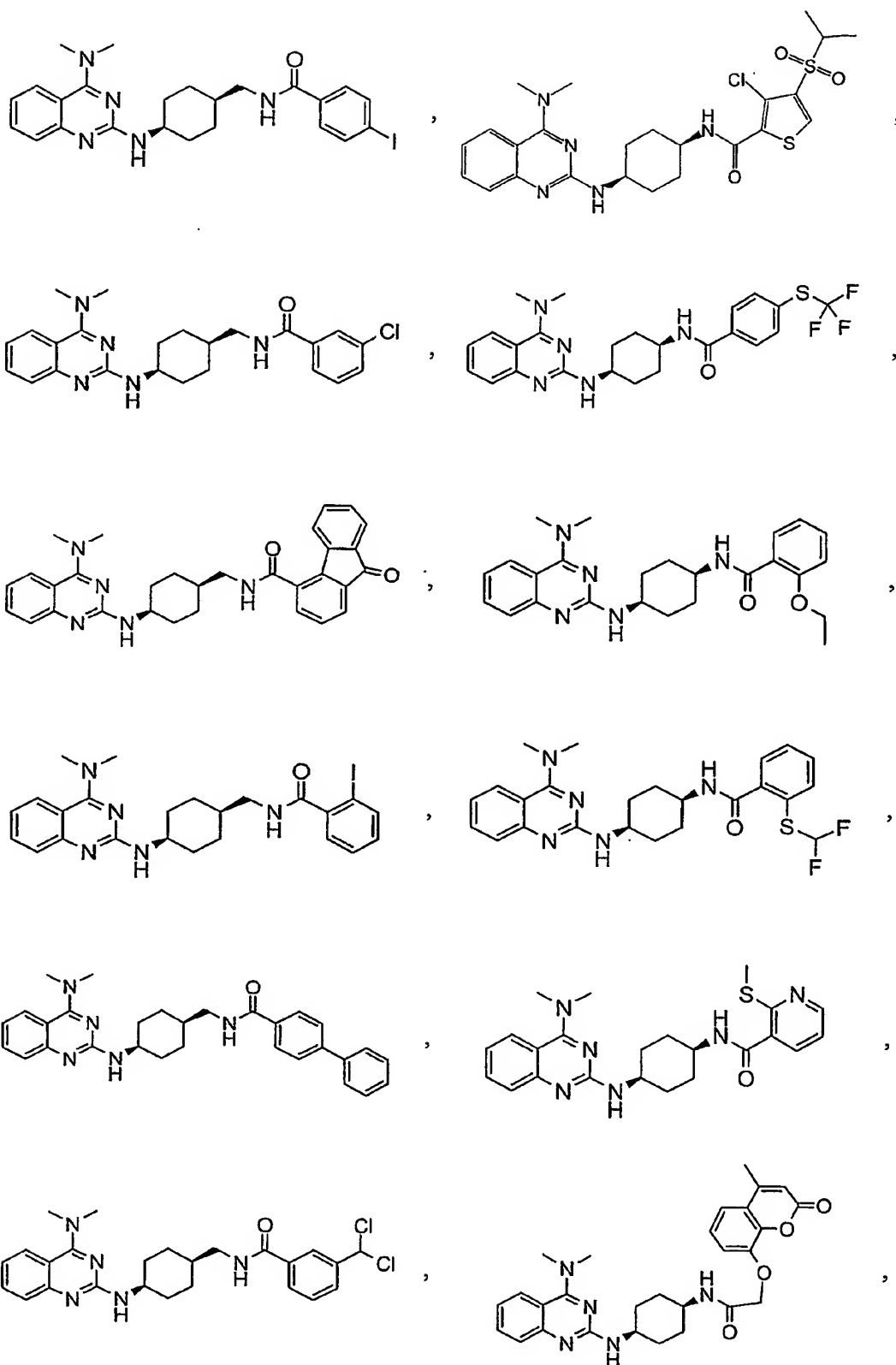


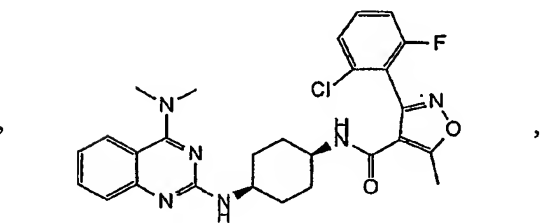
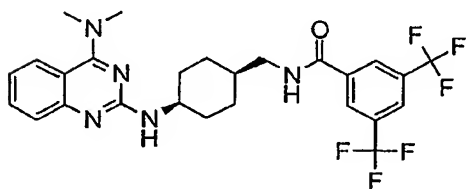
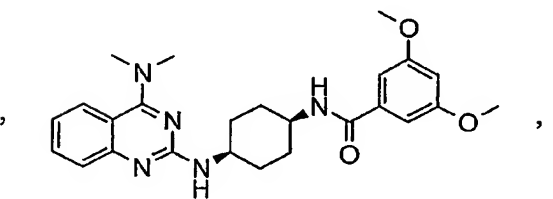
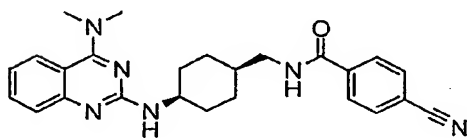
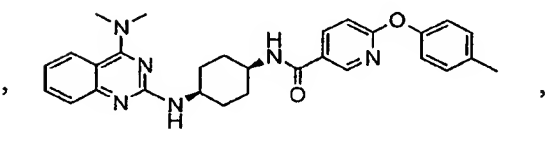
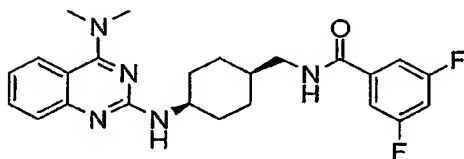
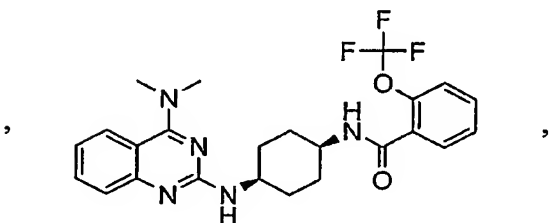
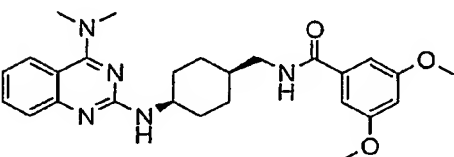
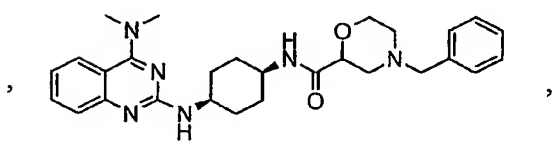
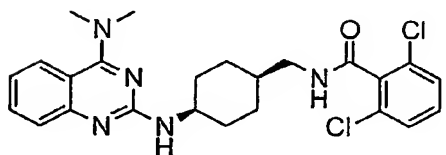
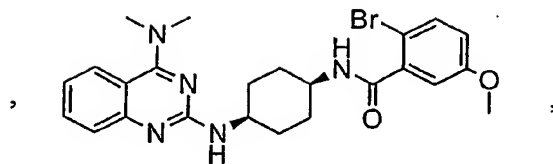
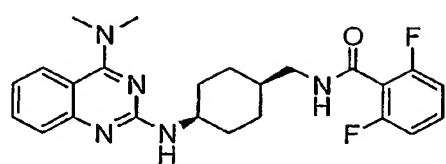


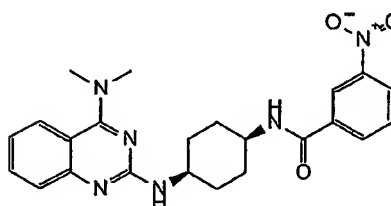
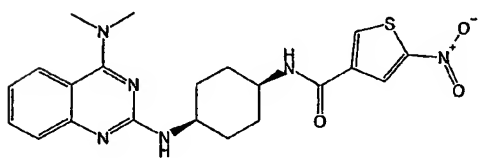
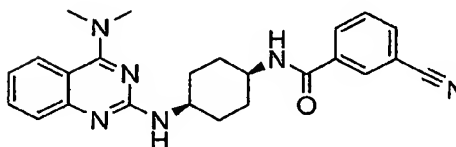
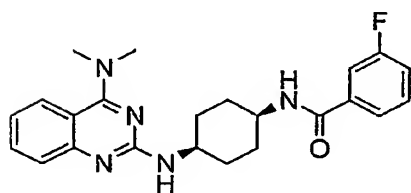
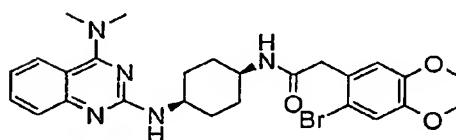
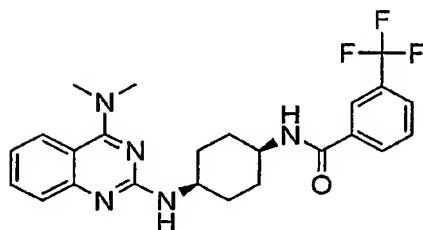
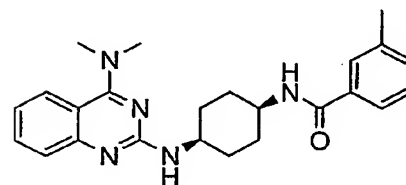
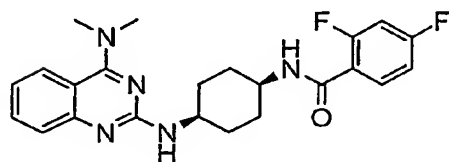
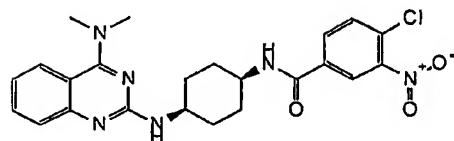
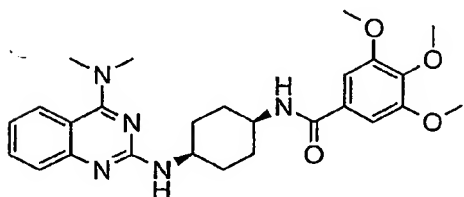
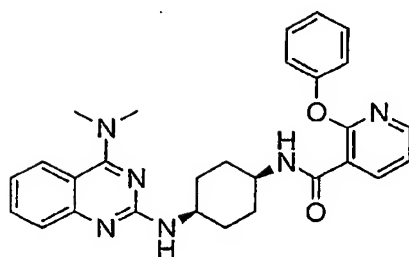
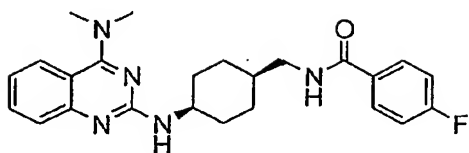


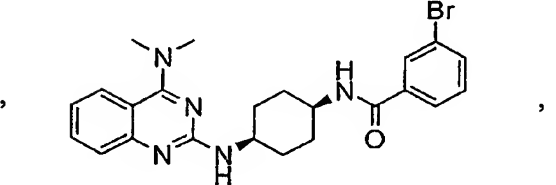
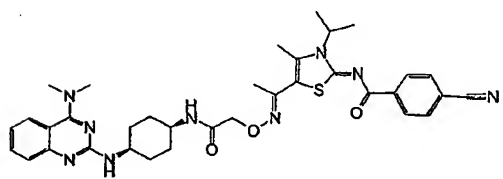
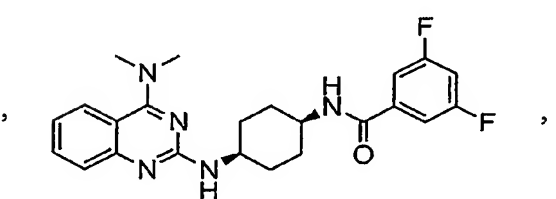
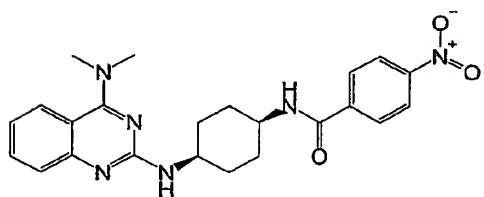
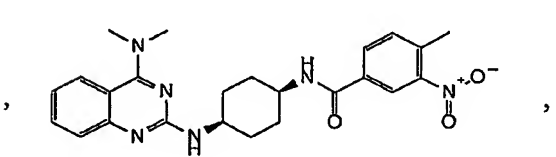
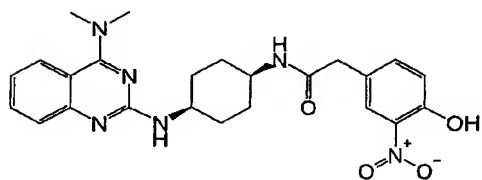
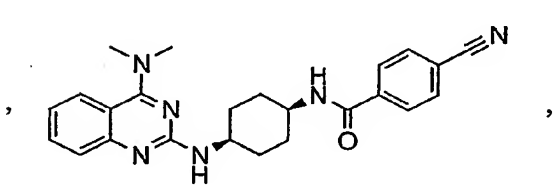
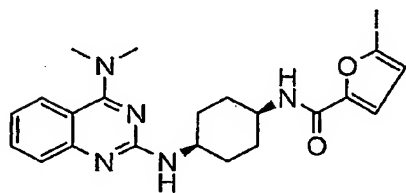
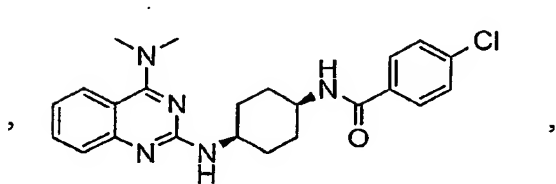
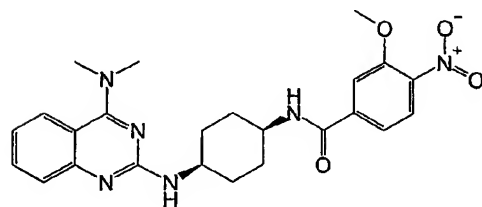
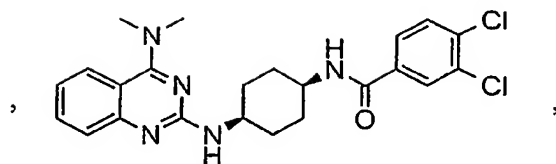
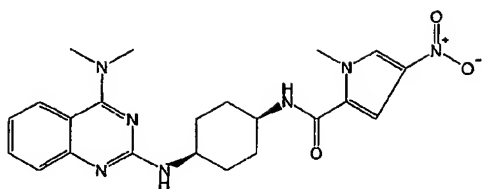


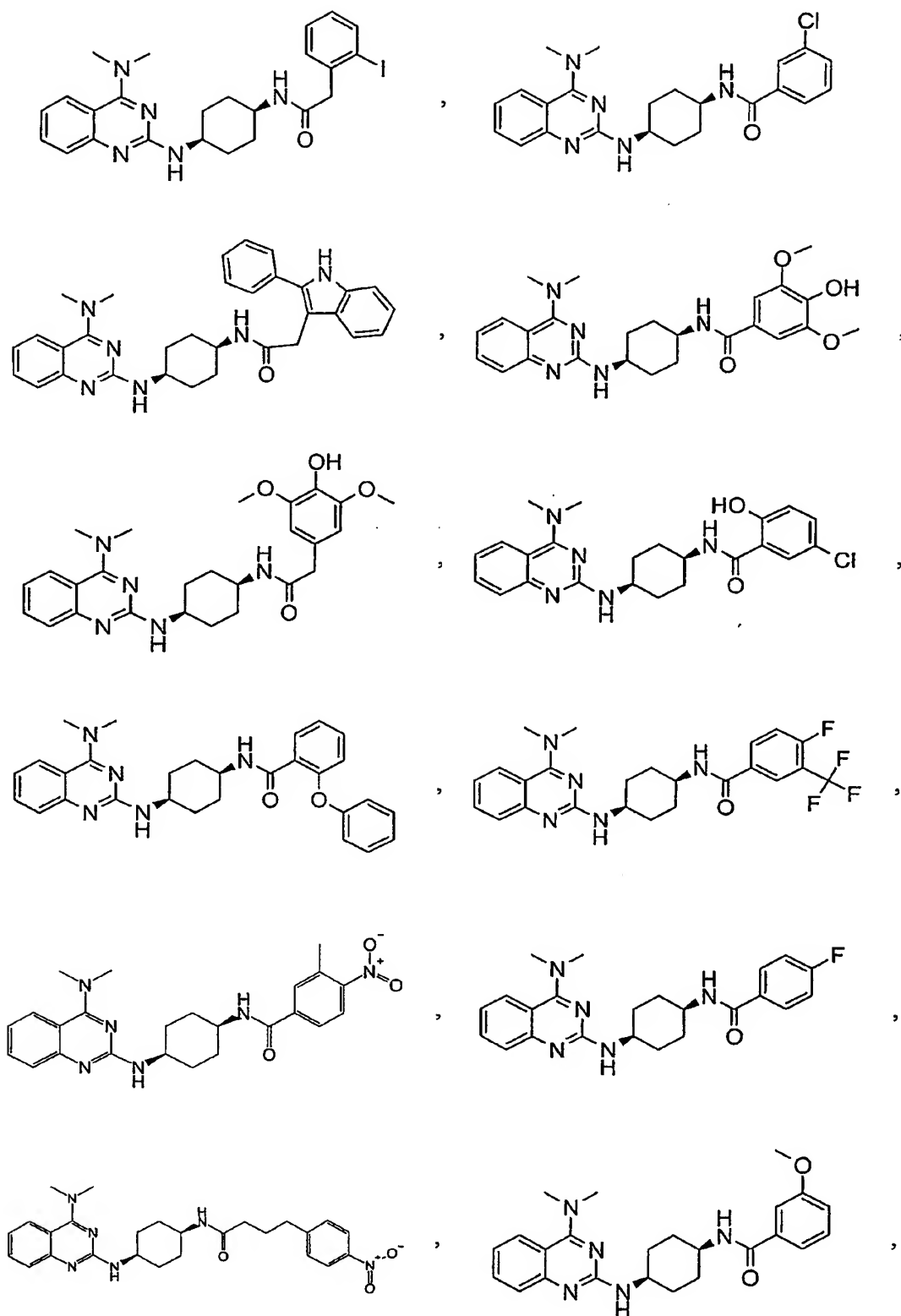


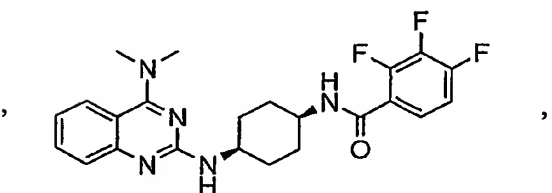
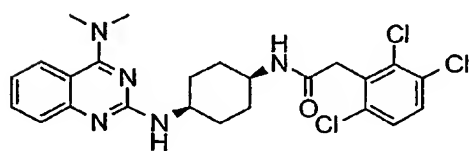
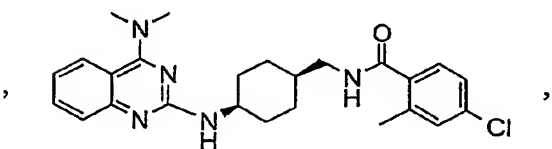
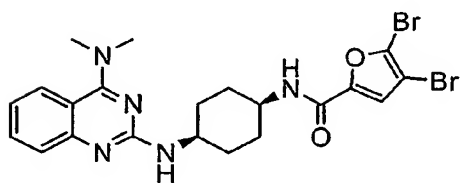
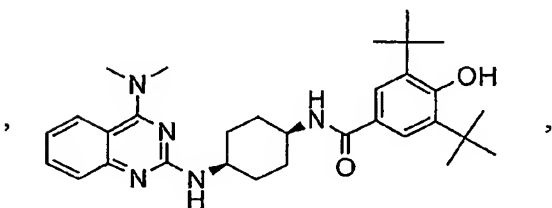
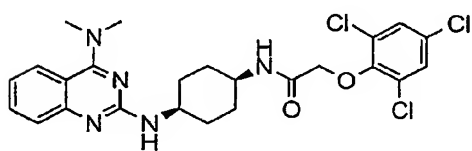
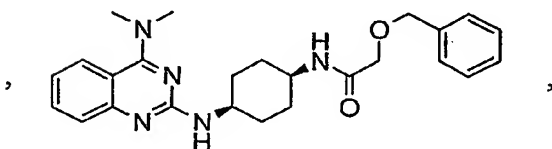
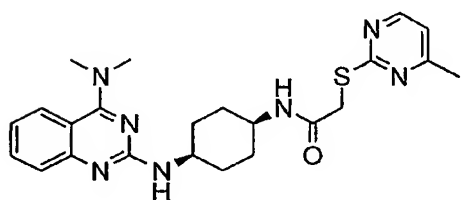
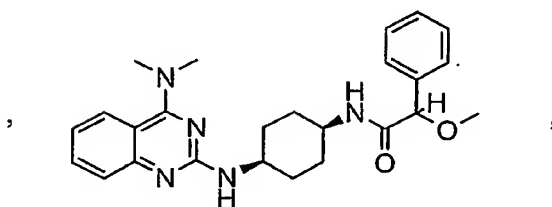
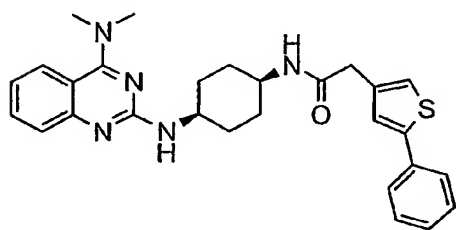
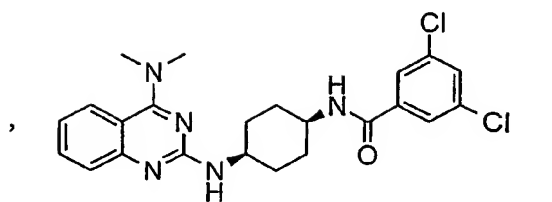
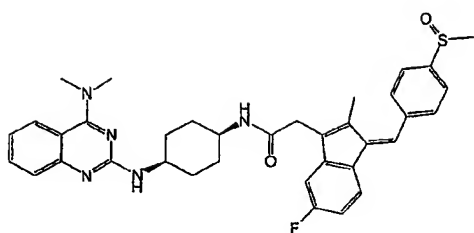


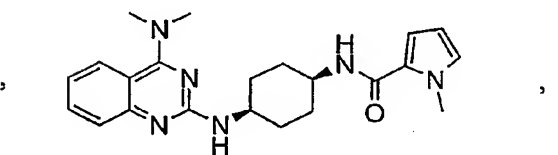
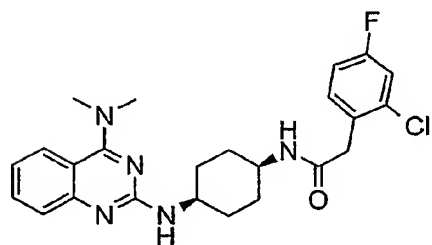
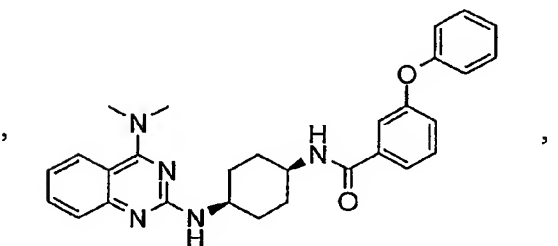
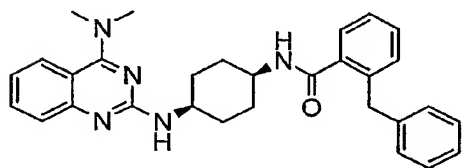
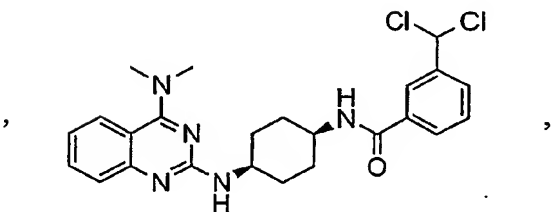
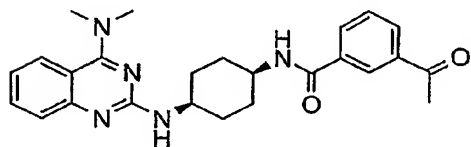
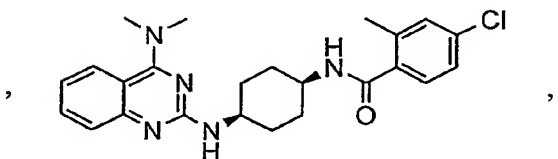
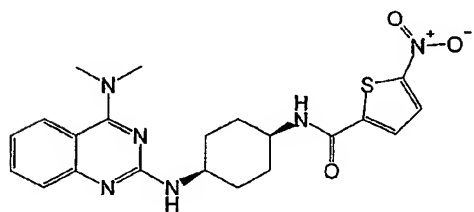
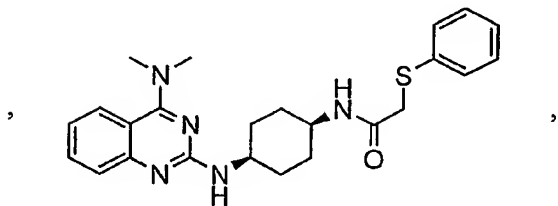
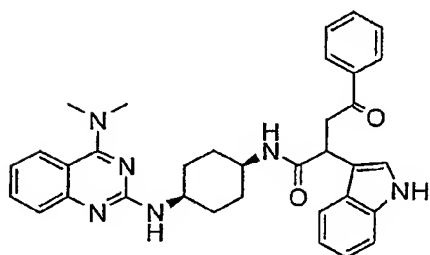
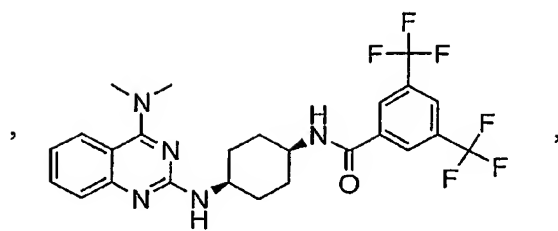
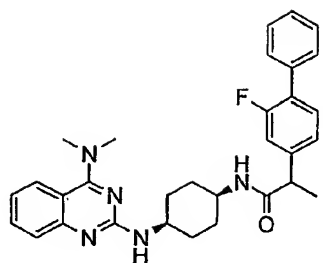


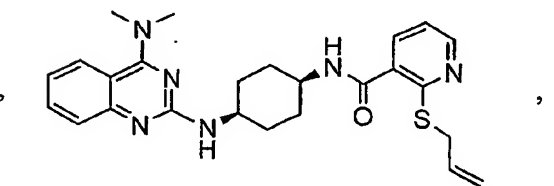
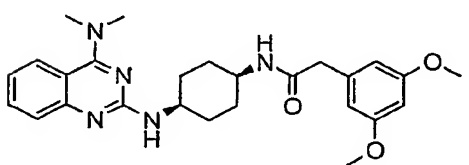
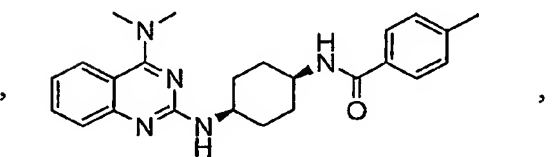
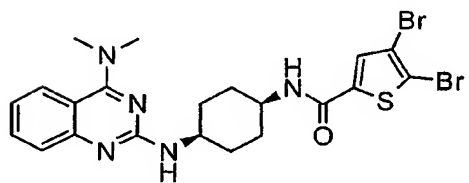
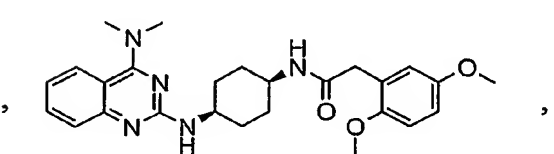
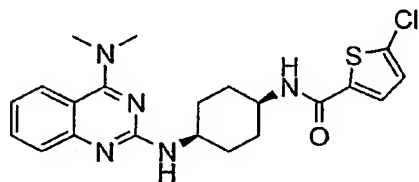
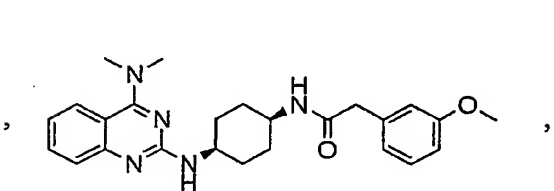
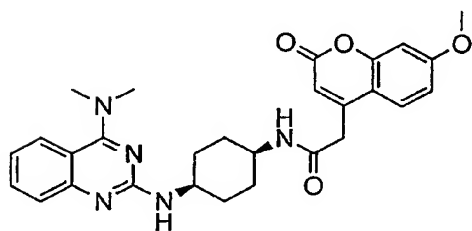
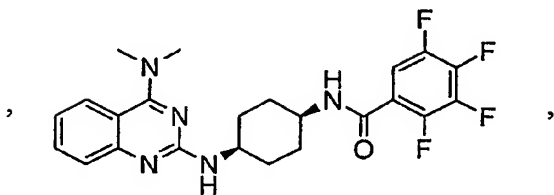
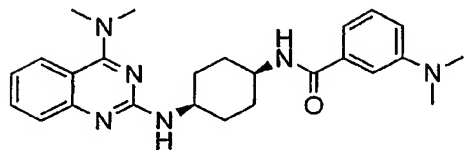
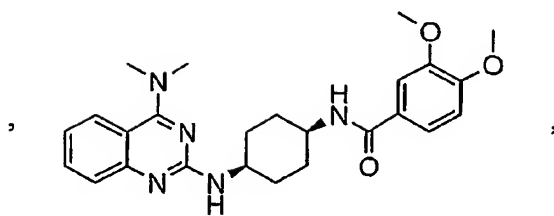
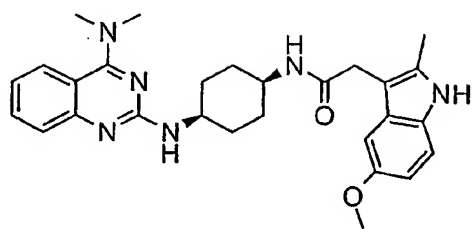


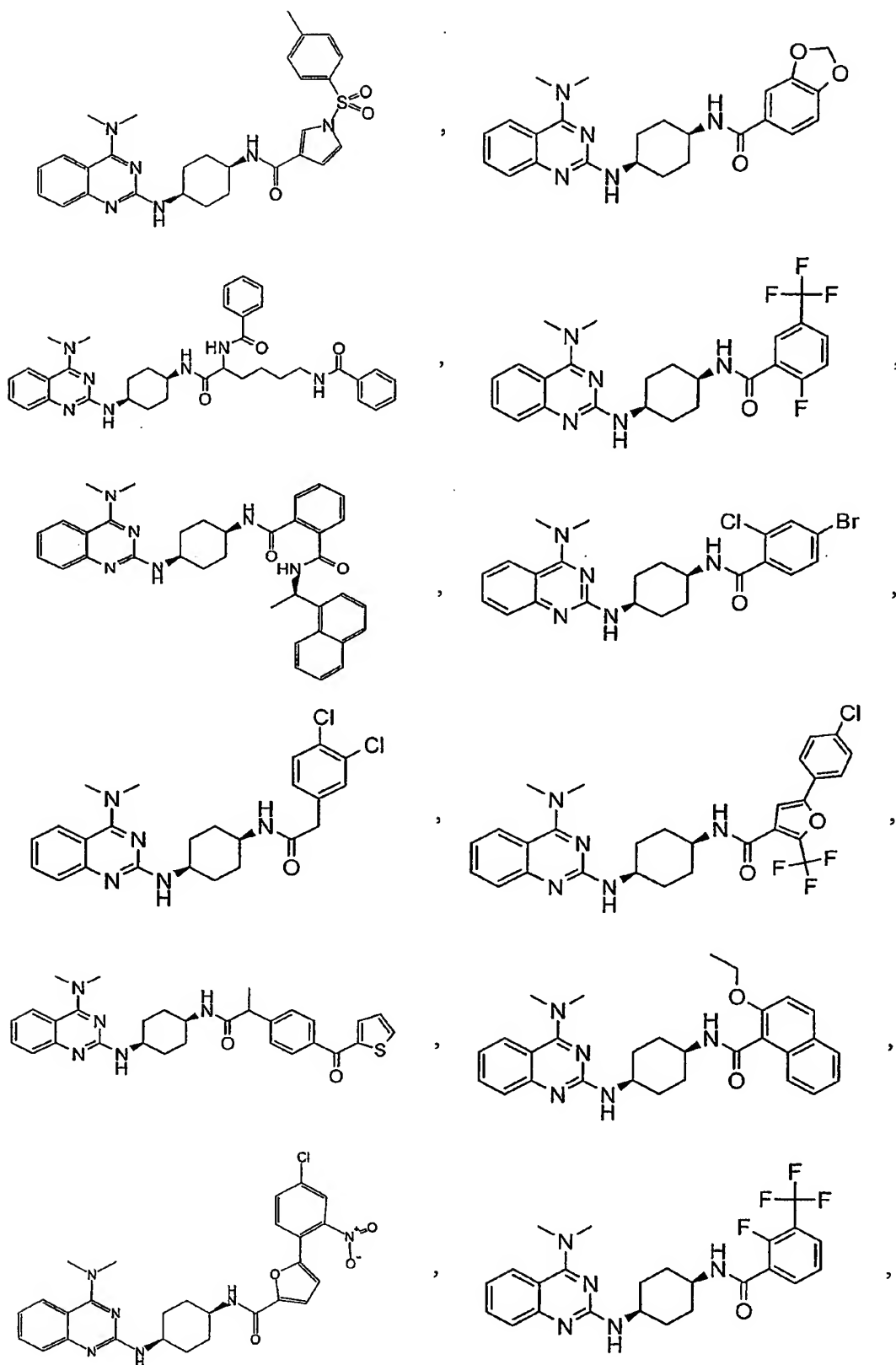


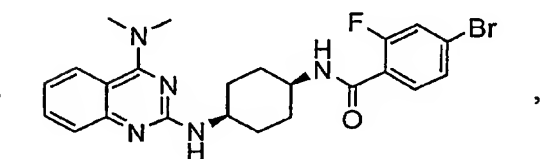
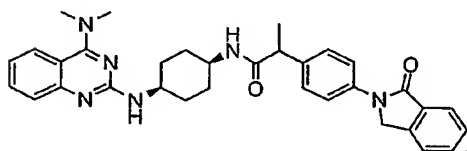
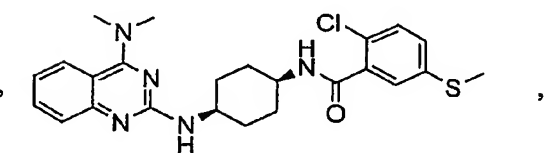
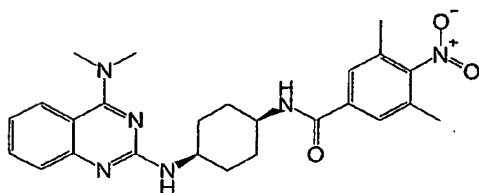
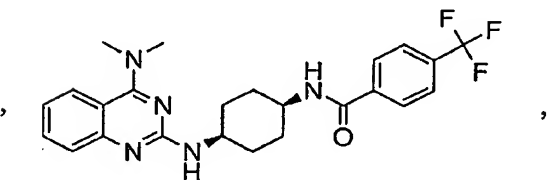
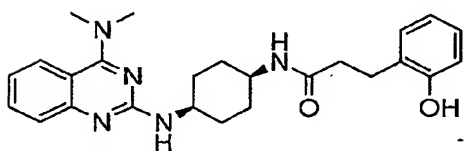
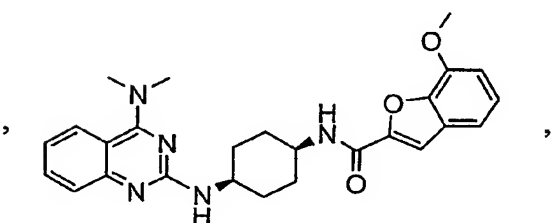
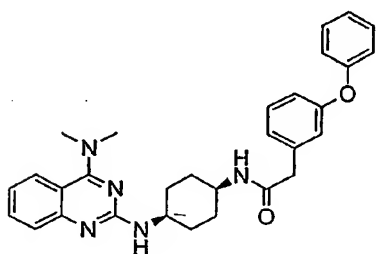
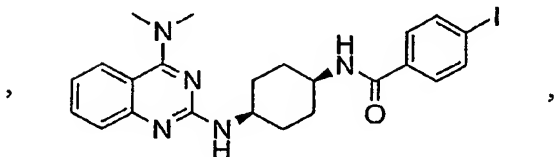
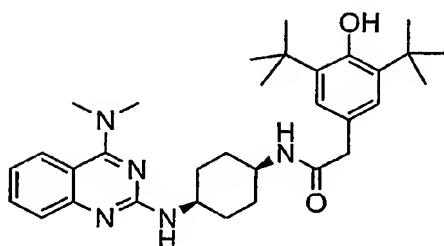
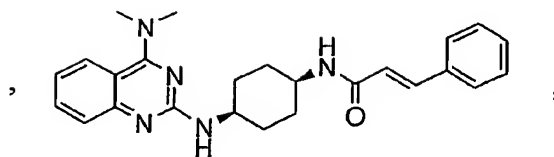
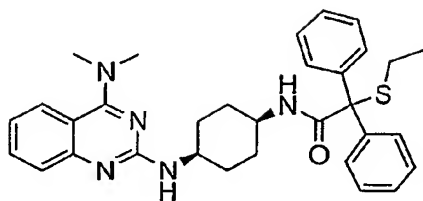


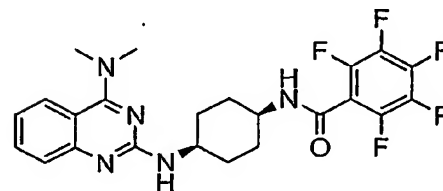
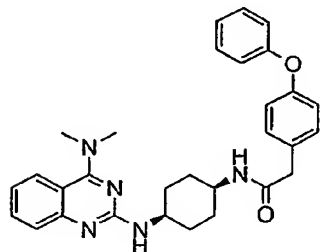
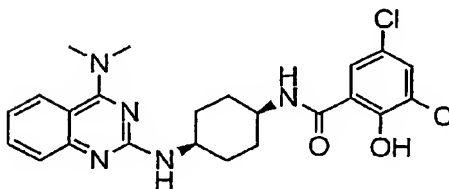
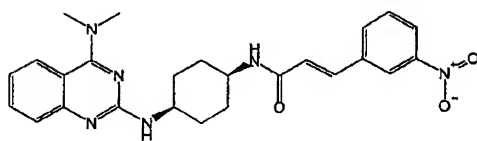
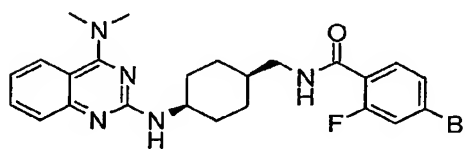
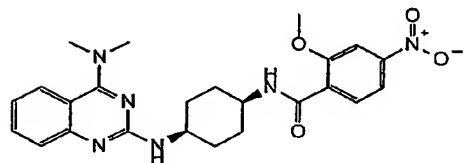
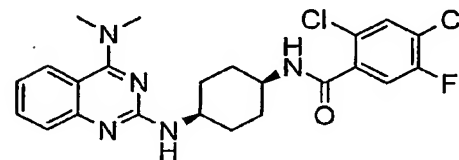
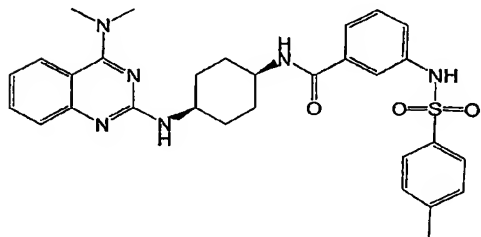
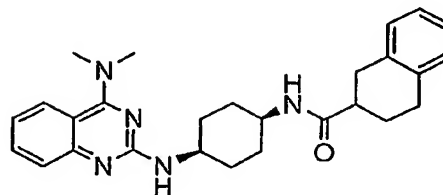
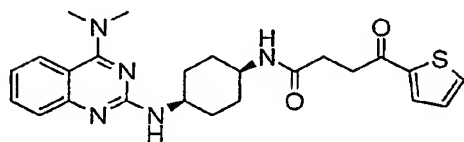
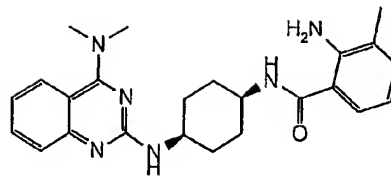
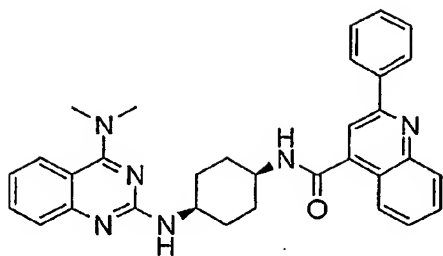


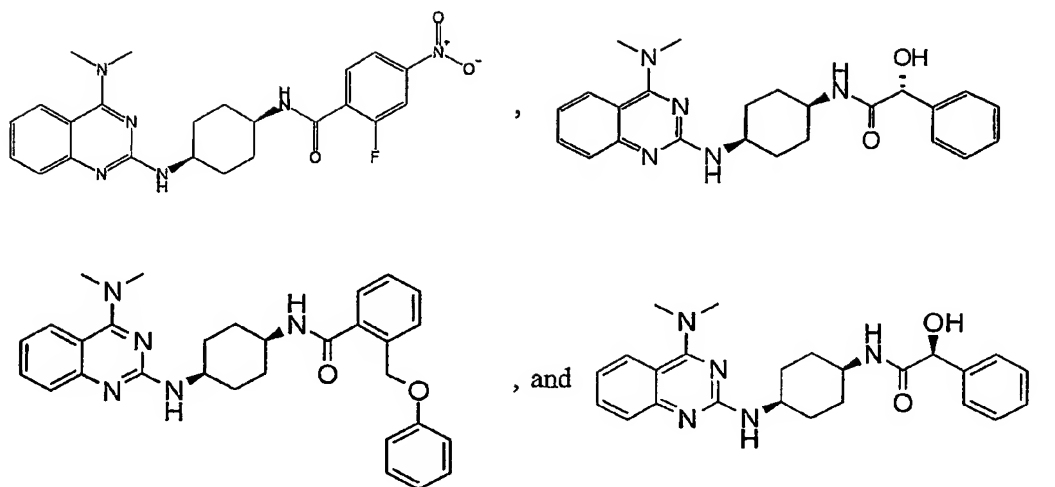












; or, in case of, a salt thereof.

7. A compound according to claim 3, wherein

R₁ represents

(i) C₁-C₁₀ alkyl,

C₁-C₁₀ alkyl substituted by substituent(s) independently selected from

•C₅-C₆ cycloalkyl,

•carbocyclic aryl,

•heterocyclyl,

(ii) C₃-C₆ cycloalkyl,

(iii) carbocyclic aryl,

(iv) or heterocyclyl;

L is selected from Formula XX - XXII;

wherein carbocyclic aryl is phenyl, naphthyl, anthranyl, or biphenyl;

heterocyclyl is 1,3-dioxo-isindolyl, 1*H*-indolyl, 1-oxo-3*H*-isobenzofuranyl, 2,3-dihydro-benzo[1,4]dioxinyl, 3,4-dihydro-2*H*-benzo[b][1,4]dioxepinyl, 4-oxo-3,4-dihydro-phthalazinyl, 9,10,10-trioxo-thioxanthenyl, 9*H*-xanthenyl, benzimidazolyl, benzo[1,3]dioxolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, furyl, imidazolyl, isoxazolyl, morpholino, oxolanyl, piperidyl, pyridyl, quinoxalyl, thienyl, quinolyl, or benzothiazolyl; or a salt thereof.

8. A compound according to claim 7, wherein

R₁ represents

(i) C₁-C₄ alkyl,

C₁-C₄ alkyl substituted by substituent(s) independently selected from

•cyclopentyl,

•carbocyclic aryl,

•heterocyclyl,

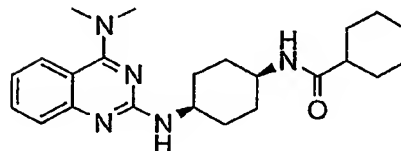
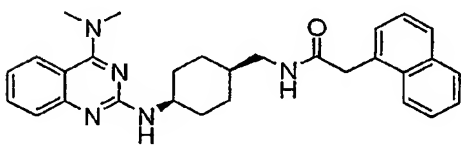
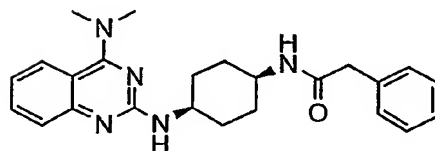
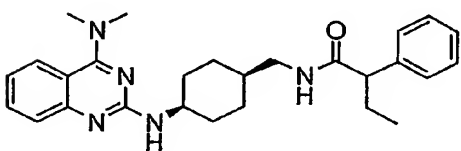
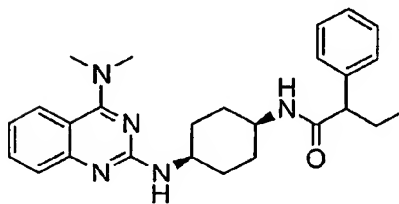
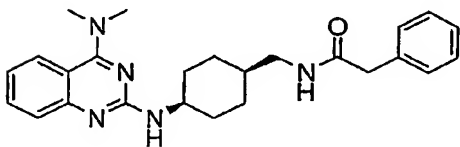
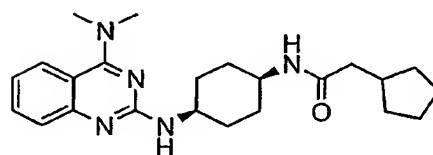
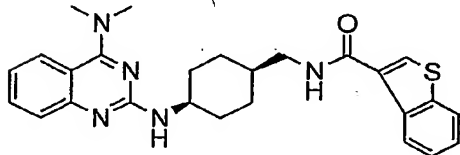
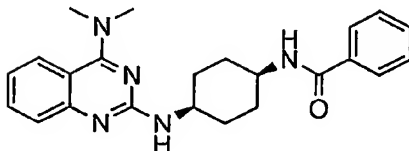
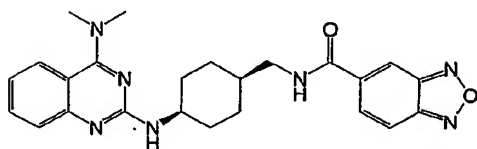
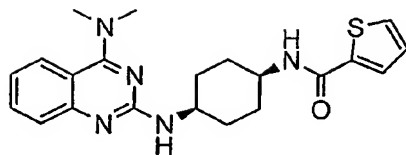
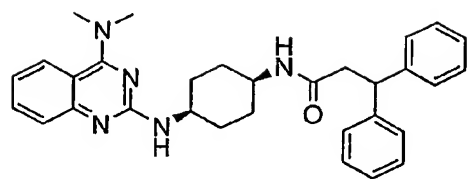
(ii) carbocyclic aryl,

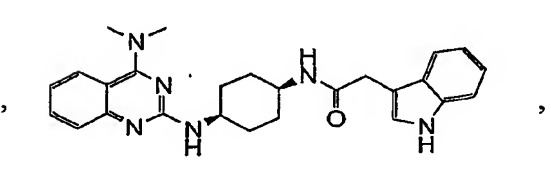
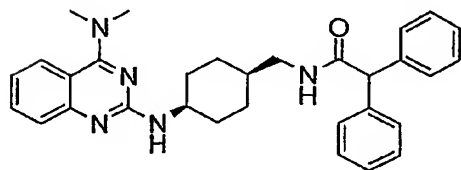
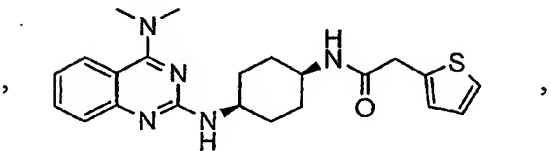
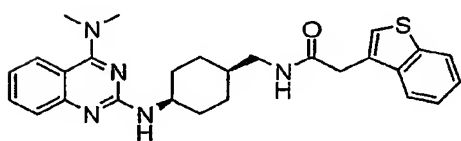
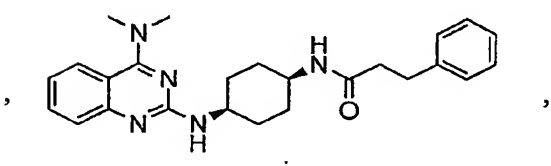
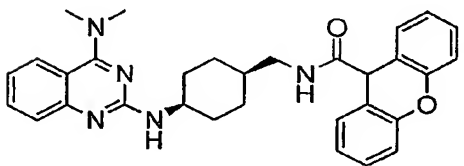
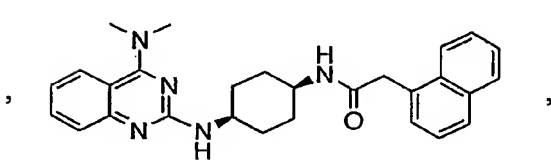
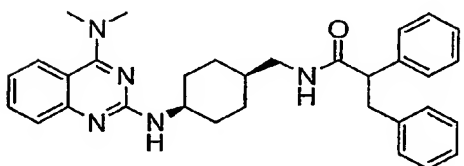
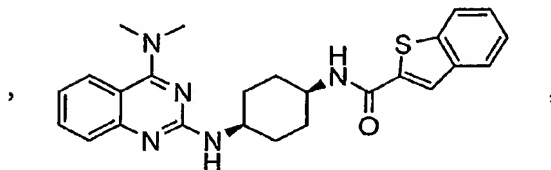
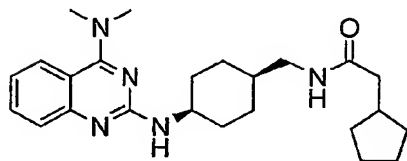
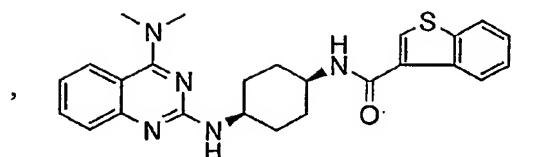
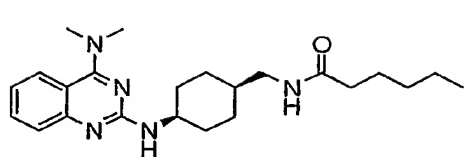
(iii) or heterocyclyl;

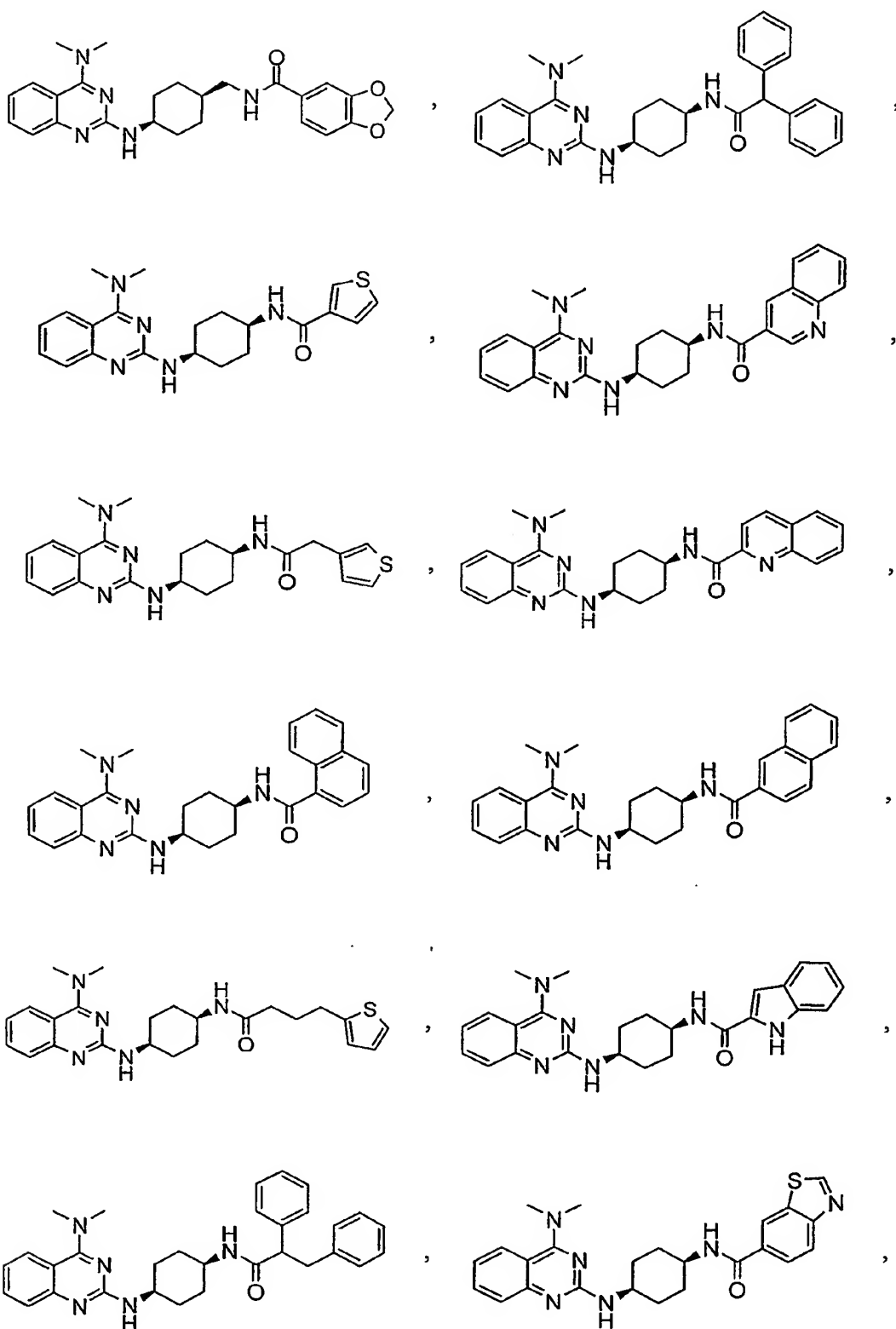
wherein carbocyclic aryl is phenyl, naphthyl, anthranyl, or biphenyl;

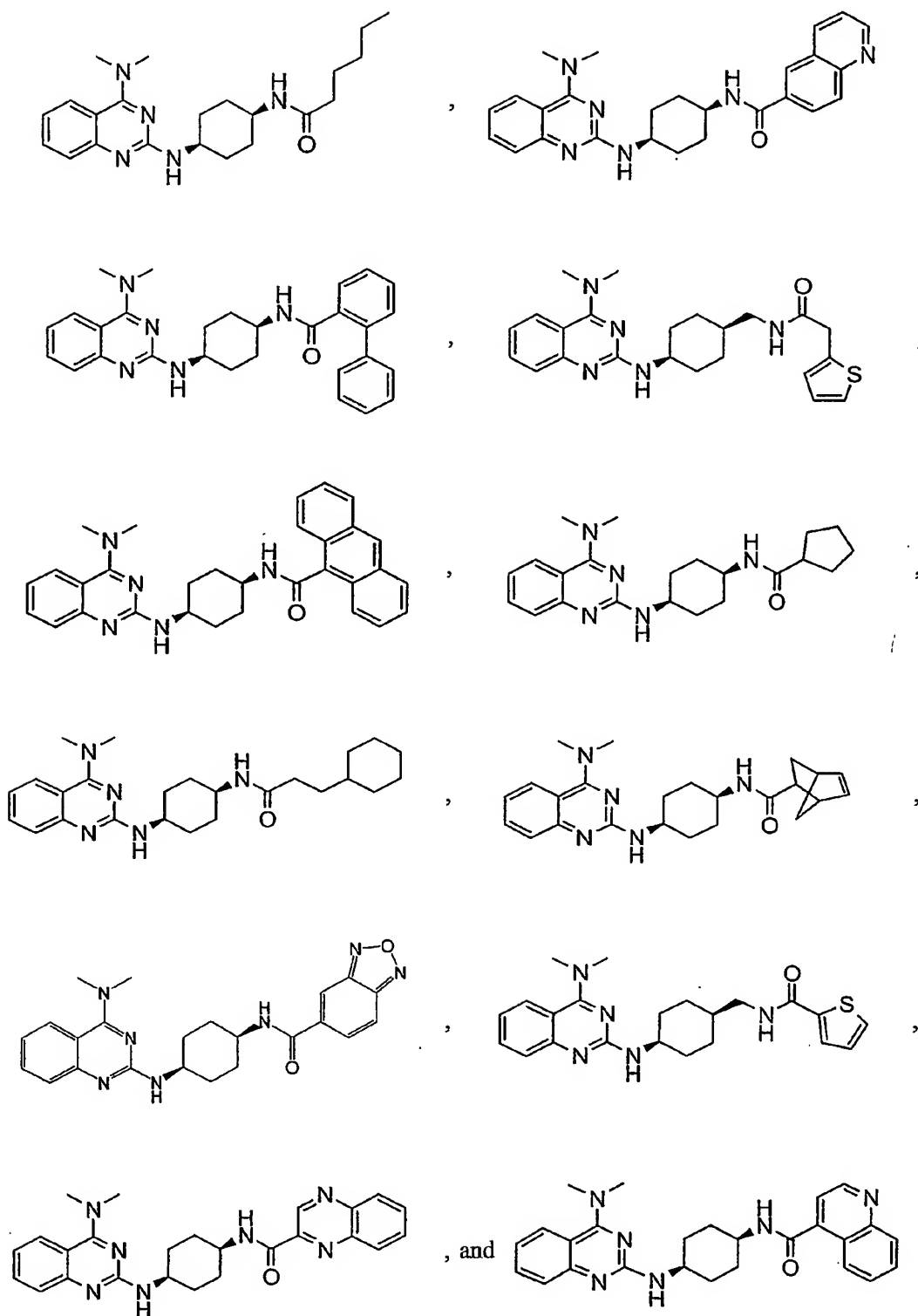
heterocyclyl is 9*H*-xanthenyl, benzo[1,3]dioxolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, thienyl, 1*H*-indolyl, quinoxalyl, quinolyl, or benzothiazolyl; or a salt thereof.

9. A compound according to claim 8 of Formua I thereof selected from the group consisting of









; or, in case of, a salt thereof.

10. A compound according to claim 1, wherein Q is Formula II;

R₁ represents

(i) C₁-C₁₀ alkyl,

C₁-C₁₀ alkyl substituted by substituent(s) independently selected from

•halogen,

•hydroxy,

•oxo,

•C₁-C₃ alkoxy,

•C₁-C₃ alkoxy substituted by substituent(s) independently selected from

••carbocyclic aryl,

••heterocyclyl,

••heterocyclyl substituted by C₁-C₃ alkyl,

•carbocyclic aryloxy,

•carbocyclic aryloxy substituted by substituent(s) independently selected from

••halogen,

••nitro,

••carbocyclic aryl,

••carbocyclic aryl substituted by C₁-C₃ alkoxy,

••C₁-C₄ alkyl,

••C₁-C₄ alkyl substituted by substituent(s) independently selected from

•••mono- or di-C₁-C₃ alkylamino,

•••mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,

•••mono- or di-C₁-C₃ alkylamino substituted by halogenated carbocyclic aryl,

•mono- or di-C₁-C₃ alkylamino,

•mono- or di-C₁-C₃ alkylamino substituted by substituent(s) independently selected from

••cyano,

••carbocyclic aryl,

••heterocyclyl,

•mono- or di-carbocyclic arylamino,

•mono- or di-carbocyclic arylamino substituted by C₁-C₃ alkyl,

•C₁-C₃ alkylcarbonylamino,

•C₁-C₄ alkoxycarbonylamino,

- carbocyclic arylsulfonylamino,
- carbocyclic arylsulfonylamino substituted by substituent(s) independently selected from
 - nitro,
 - C₁-C₃ alkyl,
 - mono- or di-C₁-C₃ alkylamino,
 - C₁-C₃ alkylthio,
 - C₁-C₃ alkylthio substituted by substituent(s) independently selected from
 - mono- or di-carbocyclic arylamino,
 - halogenated mono- or di-carbocyclic arylamino,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkoxy,
 - carbocyclic arylthio,
 - carbocyclic arylthio substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - carbocyclic arylsulfonyl,
 - halogenated carbocyclic arylsulfonyl,
 - heterocyclylthio,
 - C₃-C₆ cycloalkyl,
 - C₃-C₆ cycloalkyl substituted by C₁-C₃ alkyl,
 - carbocyclyl,
 - carbocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - C₂-C₃ alkenyl,
 - C₂-C₃ alkenyl substituted by carbocyclic aryl,
 - C₂-C₃ alkenyl substituted by carbocyclic aryl substituted C₁-C₃ alkylsulfinyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,

- hydroxy,
- nitro,
- C₁-C₄ alkyl,
- C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - carbocyclic aryl,
 - mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylamino substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by substituent(s) independently selected from
 - halogen,
 - carbocyclic aryl,
 - carbocyclic aryloxy,
 - C₁-C₃ alkoxycarbonyl,
 - mono- or di-C₁-C₃ alkylamino,
 - C₁-C₃ alkylthio,
 - halogenated C₁-C₃ alkylthio,
 - C₁-C₃ alkylsulfonyl,
 - C₃-C₆ cycloalkyl,
 - carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - C₁-C₃ alkoxy,

- C₁-C₃ alkoxy substituted by carbocyclic aryl,
- carbocyclic aryl,
- halogenated carbocyclic aryl,
- (ii) C₂-C₈ alkenyl,
- C₂-C₈ alkenyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - C₁-C₃ alkoxy,
 - halogenated C₁-C₃ alkoxy,
 - heterocyclyl,
 - heterocyclyl substituted by nitro,
- (iii) C₂-C₄ alkynyl,
- C₂-C₄ alkynyl substituted by carbocyclic aryl,
- (iv) C₃-C₆ cycloalkyl,
- C₃-C₆ cycloalkyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - C₁-C₃ alkyl substituted by substituent(s) independently selected from
 - hydroxy,
 - oxo,
 - carbocyclic aryl,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,
 - carbocyclic aryl,
- (v) C₃-C₆ cycloalkenyl,
- C₃-C₆ cycloalkenyl substituted by C₁-C₃ alkyl,
- (vi) carbocyclyl,
- carbocyclyl substituted by substituent(s) independently selected from

- hydroxy,
- nitro,
- (vii) carbocyclic aryl,
carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₉ alkyl,
 - C₁-C₉ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - oxo,
 - C₁-C₃ alkoxy,
 - carbocyclic aryloxy,
 - mono- or di-C₁-C₃ alkylamino-N-oxy,
 - mono- or di-C₁-C₃ alkylamino,
 - mono- or di-C₁-C₃ alkylamino substituted by carbocyclic aryl,
 - mono- or di-carbocyclic arylamino,
 - mono- or di-carbocyclic arylamino substituted by C₁-C₃ alkoxy,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
 - heterocyclyl,
 - heterocyclyl substituted by C₁-C₃ alkyl,
 - C₂-C₃ alkenyl,
 - C₂-C₃ alkenyl substituted by carbocyclic aryl,
 - C₁-C₉ alkoxy,
 - C₁-C₉ alkoxy substituted by substituent(s) independently selected from
 - hydroxy,
 - halogen,
 - carboxy,
 - mono- or di-C₁-C₃ alkylamino,

- carbocyclic aryl,
- halogenated carbocyclic aryl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
- halogen,
- C₁-C₃ alkyl,
- halogenated C₁-C₃ alkyl,
- C₂-C₃ alkenyloxy,
- C₁-C₃ alkylcarbonyloxy,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by substituent(s) independently selected from
- halogen,
- C₁-C₄ alkyl,
- halogenated C₁-C₄ alkyl,
- C₁-C₃ alkoxy,
- heterocyclyloxy,
- heterocyclyloxy substituted by substituent(s) independently selected from
- halogen,
- C₁-C₃ alkyl,
- halogenated C₁-C₃ alkyl,
- (carbocyclic aryl)S(O)₂O,
- carboxy,
- C₁-C₃ alkoxycarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl,
- mono- or di-C₁-C₃ alkylaminocarbonyl substituted by carbocyclic aryl,
- amino,
- mono- or di-C₁-C₄ alkylamino,
- mono- or di-C₁-C₄ alkylamino substituted by cyano,
- mono- or di-carbocyclic arylamino,
- C₁-C₃ alkylcarbonylamino,

- carbocyclic arylsulfonylamino,
- carbocyclic arylsulfonylamino substituted by C₁-C₃ alkyl,
- (carbocyclic aryl)NHC(O)NH,
- (carbocyclic aryl)NHC(O)NH substituted by C₁-C₃ alkoxy,
- (carbocyclic aryl)NHC(O)NH substituted by halogenated C₁-C₃ alkoxy,
- C₁-C₃ alkylthio,
- halogenated C₁-C₃ alkylthio,
- carbocyclic arylthio,
- halogenated carbocyclic arylthio,
- carbocyclic arylthio substituted by C₁-C₃ alkyl,
- heterocyclylthio,
- C₁-C₃ alkylsulfonyl,
- mono- or di-C₁-C₃ alkylaminosulfonyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - C₁-C₇ alkyl,
 - halogenated C₁-C₇ alkyl,
 - heterocyclyl,
 - heterocyclyl substituted by substituent(s) independently selected from
 - C₁-C₃ alkyl,
 - carbocyclic aryl,
 - halogenated carbocyclic aryl,
- (viii) heterocyclyl,
- or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,

- oxo,
- C₁-C₃ alkylcarbonyloxy,
- C₁-C₃ alkoxycarbonyl,
- C₁-C₃ alkylthio,
- C₁-C₃ alkylthio substituted by carbocyclic aryl,
- C₁-C₃ alkylthio substituted by halogenated carbocyclic aryl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - heterocyclyl,
- C₁-C₃ alkoxy,
- C₁-C₃ alkoxy substituted by carbocyclic aryl,
- carbocyclic aryloxy,
- carbocyclic aryloxy substituted by C₁-C₃ alkyl,
- mono- or di-C₁-C₃ alkylamino,
- C₁-C₄ alkylcarbonylamino,
- C₁-C₃ alkylthio,
- carbocyclic arylthio,
- halogenated carbocyclic arylthio,
- carbocyclic arylthio substituted by C₁-C₃ alkoxycarbonyl,
- heterocyclylthio,
- heterocyclylthio substituted by C₁-C₃ alkyl,
- C₁-C₃ alkylsulfonyl,
- carbocyclic arylsulfonyl,
- carbocyclic arylsulfonyl substituted by C₁-C₄ alkyl,
- C₁-C₃ alkoxycarbonyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₃ alkyl,

- halogenated C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- halogenated C₁-C₃ alkoxy,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
- C₁-C₃ alkyl,
- halogenated C₁-C₃ alkyl,
- C₁-C₃ alkoxy,
- C₁-C₃ alkoxycarbonyl;

Y is -(CH₂)_m, m is 0 or 1;

wherein carbocyclic aryl is phenyl, naphthyl, biphenyl, or phenanthryl;

carbocyclyl is 9*H*-fluorenyl, 9-oxo-fluorenyl, acenaphthyl, anthraquinonyl, indanyl, or indenyl;

heterocyclyl is 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1,3,4-thiadiazolyl, 1,3-dioxo-isoindolyl, 1,3-dioxolanyl, 1*H*-indolyl, 1*H*-pyrrolo[2,3-*c*]pyridyl, 1*H*-pyrrolyl, 2,2',5',2''-terthiophenyl, 2,2'-bithiophenyl, 2,3-dihydro-1-oxo-isoindolyl, 2,3-dihydro-benzo[1,4]dioxinyl, 2,3-dihydro-benzofuryl, 2,4-dihydro-3-oxo-pyrazolyl, 2*H*-benzopyranyl, 2-oxo-pyrrolidinyl, 3,4-dihydro-2*H*-benzo[1,4]oxazinyl, 3,4-dihydro-2*H*-benzo[*b*][1,4]dioxepinyl, 4*H*-benzo[1,3]dioxinyl, 4*H*-benzopyranyl, 4-oxo-1,5,6,7-tetrahydro-indolyl, 4-oxo-benzopyranyl, 9*H*-carbazolyl, 9*H*-xanthenyl, azetidiny, benzimidazolyl, benzo[1,3]dioxolyl, benzo[*b*]thienyl, benzofuryl, benzothiazolyl, furyl, imidazo[2,1-*b*]thiazolyl, imidazolyl, isoxazolyl, morpholino, morpholinyl, oxolanyl, piperazyl, piperidyl, pyrazolo[5,1-*b*]thiazolyl, pyrazolyl, pyridyl, pyrimidyl, pyrrolidyl, quinolyl, quinoxalyl, thiazolidyl, thiazolyl, thienyl, or thiolanyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

11. A compound according to claim 10, wherein

R₁ represents

- (i) C₁-C₁₀ alkyl substituted by substituent(s) independently selected from
- methoxy,
- methoxy substituted by carbocyclic aryl,

- carbocyclic aryloxy,
- halogenated carbocyclic aryloxy,
- mono-C₁-C₂ alkylamino substituted by cyano,
- mono- or di-C₁-C₂ alkylamino substituted by carbocyclic aryl,
- mono-carbocyclic arylamino,
- mono-carbocyclic arylamino substituted by methyl,
- carbocyclic arylsulfonylamino substituted by methyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by carbocyclic aryl,
 - C₁-C₄ alkyl substituted by hydroxy,
 - C₁-C₂ alkoxy,
 - halogenated C₁-C₂ alkoxy,
 - heterocyclyl substituted by carbocyclic aryl,
- (ii) C₂-C₈ alkenyl substituted by substituent(s) independently selected from
 - methoxy substituted by carbocyclic aryl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by methoxy,
- (iii) C₂-C₄ alkynyl substituted by carbocyclic aryl,
- (iv) cyclohexyl substituted by carbocyclic arylmethyl,
- (v) carbocyclyl,
- (vi) carbocyclic aryl,
carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - amino,
 - C₁-C₉ alkyl,
 - halogenated C₁-C₉ alkyl,

- C₁-C₉ alkoxy,
- C₁-C₉ alkoxy substituted by substituent(s) independently selected from
 - halogen,
 - halogenated carbocyclic aryl,
- propenyloxy,
- methylamino,
- di-C₁-C₂ alkylamino,
- di-C₁-C₂ alkylamino substituted by cyano,
- methylthio,
- halogenated methylthio,
- (vii) heterocyclyl,
- or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by hydroxy,
 - C₁-C₄ alkyl substituted by carbocyclic aryl,
 - methoxy,
 - C₁-C₂ alkoxycarbonyl,
 - carbocyclic arylthio substituted by methoxycarbonyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - halogenated methyl,
 - heterocyclyl;

R₂ is methylamino or dimethylamino;

L is selected from Formula Va, VIIIa, or IXa;

wherein carbocyclic aryl is phenyl, naphthyl, biphenyl, or phenanthryl;

carbocyclyl is 9*H*-fluorenyl, acenaphthyl, or anthraquinonyl;

heterocyclyl is 1,2,3-thiadiazolyl, 1,2,3-triazolyl, 1,2-dihydro-3-oxo-pyrazolyl, 1,3-dioxolanyl, 1*H*-indolyl, 1*H*-pyrrolyl, 2,2',5',2''-terthiophenyl, 2,2'-bithiophenyl, 2,3-dihydro-benzo[1,4]dioxinyl, 3,4-dihydro-2*H*-benzo[1,4]oxazinyl, 4-oxo-benzopyranyl, 9*H*-carbazolyl, 9*H*-xanthenyl, benzimidazolyl, benzo[1,3]dioxolyl, benzo[b]thienyl, benzofuryl,

benzothiazolyl, furyl, imidazolyl, isoxazolyl, oxolanyl, pyrazolo[5,1-b]thiazolyl, pyrazolyl, pyridyl, pyrimidyl, quinolyl, quinoxalyl, thiazolidyl, thiazolyl, thienyl, 2*H*-benzopyranyl, 4*H*-benzo[1,3]dioxinyl, azetidiny, imidazo[2,1-b]thiazolyl, morpholinyl, or 2,3-dihydro-benzofuryl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

12. A compound according to claim 11, wherein

R₁ represents

(i) C₁-C₇ alkyl substituted by substituent(s) independently selected from

- methoxy,
- methoxy substituted by carbocyclic aryl,
- carbocyclic aryloxy,
- halogenated carbocyclic aryloxy,
- mono-ethylamino substituted by cyano,
- di-methylamino substituted by carbocyclic aryl,
- mono-carbocyclic arylamino,
- mono-carbocyclic arylamino substituted by methyl,
- carbocyclic arylsulfonylamino substituted by methyl,
- carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - nitro,
 - C₁-C₄ alkyl,
 - C₁-C₄ alkyl substituted by carbocyclic aryl,
 - C₁-C₄ alkyl substituted by hydroxy,
 - methoxy,
 - halogenated methoxy,
 - heterocyclyl substituted by carbocyclic aryl,

(ii) C₂-C₇ alkenyl substituted by substituent(s) independently selected from

- methoxy substituted by carbocyclic aryl,
- carbocyclic aryl,

- carbocyclic aryl substituted by methoxy,
- (iii) butynyl substituted by carbocyclic aryl,
- (iv) cyclohexyl substituted by carbocyclic arylmethyl,
- (v) carbocyclyl,
- (vi) carbocyclic aryl,
- carbocyclic aryl substituted by substituent(s) independently selected from
 - halogen,
 - hydroxy,
 - cyano,
 - amino,
 - C₁-C₂ alkyl,
 - halogenated methyl,
 - C₁-C₃ alkoxy,
 - C₁-C₃ alkoxy substituted by substituent(s) independently selected from
 - halogen,
 - halogenated carbocyclic aryl,
 - propenyloxy,
 - di-C₁-C₂ alkylamino,
 - di-C₁-C₂ alkylamino substituted by cyano,
 - methylthio,
 - halogenated methylthio,
 - (vii) heterocyclyl,
 - or heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - C₁-C₃ alkyl substituted by hydroxy,
 - C₁-C₃ alkyl substituted by carbocyclic aryl,
 - methoxy,
 - ethoxycarbonyl,
 - carbocyclic arylthio substituted by methoxycarbonyl,
 - carbocyclic aryl,
 - carbocyclic aryl substituted by substituent(s) independently selected from

- halogen,
- halogenated methyl,
- heterocyclyl;

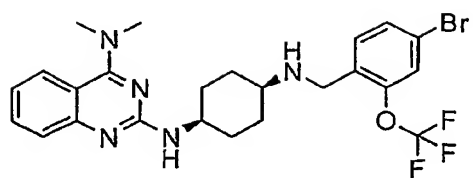
L is selected from Formula XX - XXII;

wherein carbocyclic aryl is phenyl, naphthyl, or biphenyl;

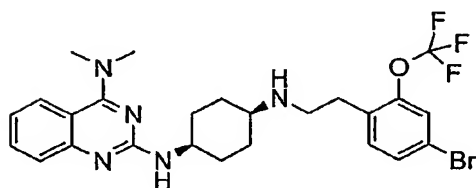
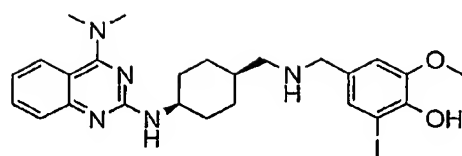
carbocyclyl is acenaphthyl;

heterocyclyl is 1*H*-indolyl, 1*H*-pyrrolyl, 2,3-dihydro-benzo[1,4]dioxinyl, 9*H*-carbazolyl, benzo[1,3]dioxolyl, furyl, pyrazolyl, thienyl, 4-oxo-benzopyranyl, azetidiny, imidazo[2,1-*b*]thiazolyl, pyridyl, imidazolyl, 2,3-dihydro-benzofuryl, or benzo[*b*]thienyl;
halogen is fluoro, chloro, bromo, or iodo;
or a salt thereof.

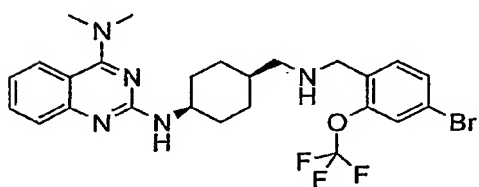
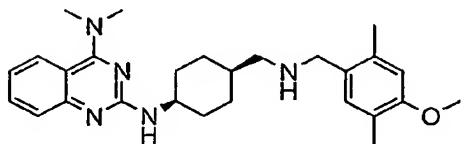
13. A compound according to claim 12 of Formula I selected from the group consisting of



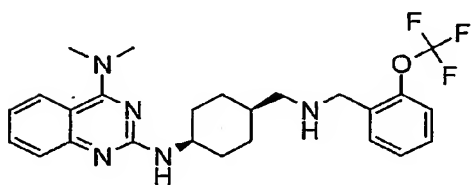
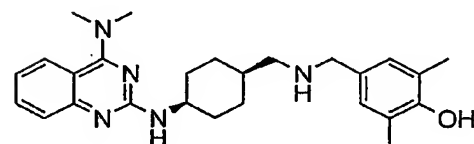
,



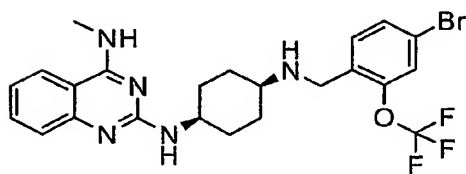
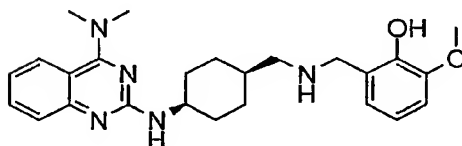
,



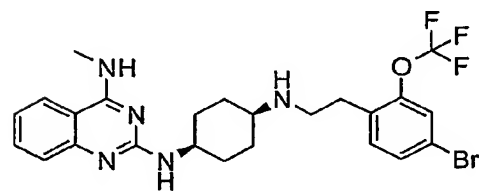
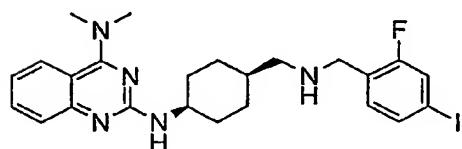
,



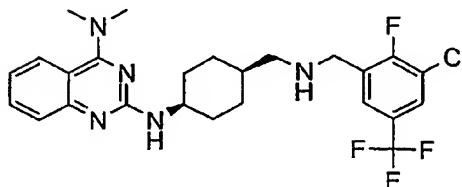
,

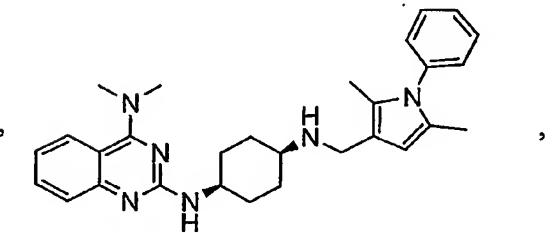
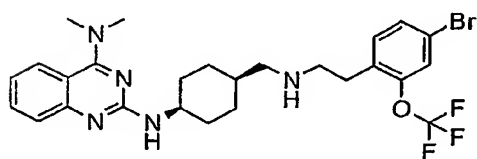
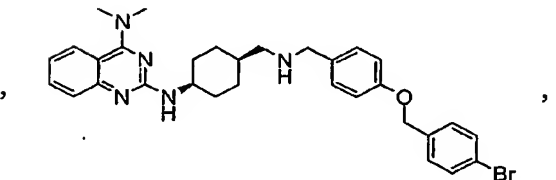
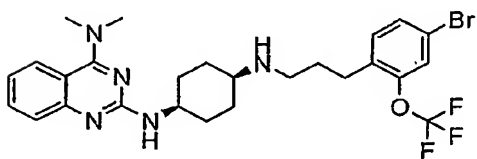
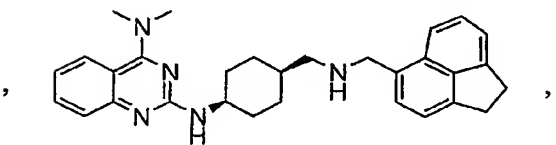
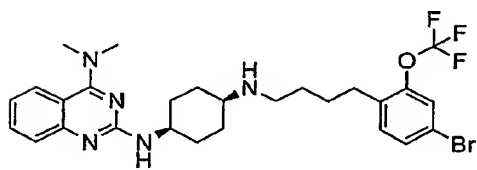
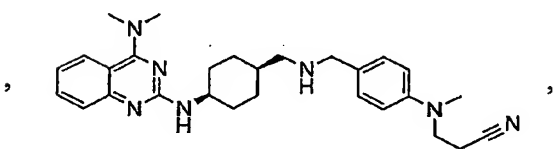
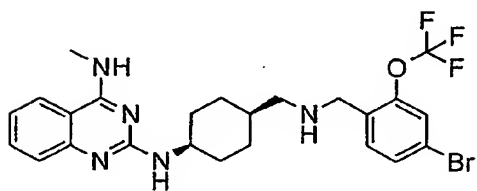
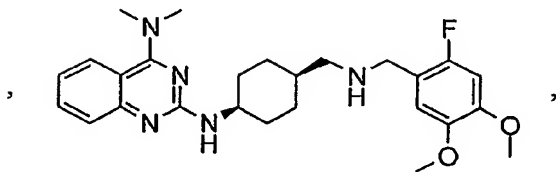
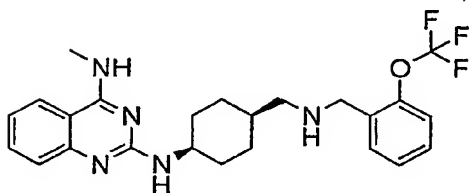
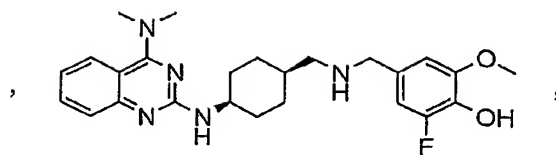
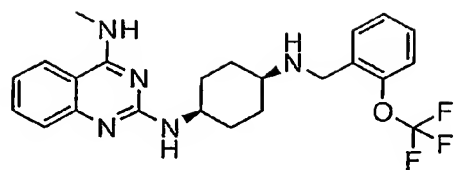


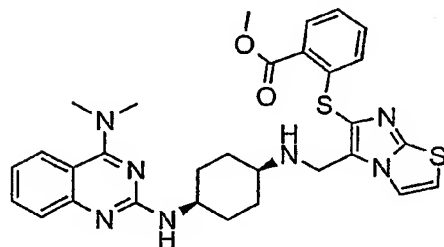
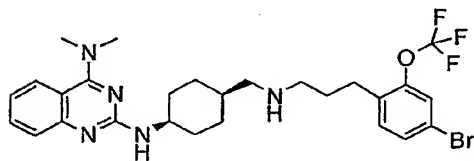
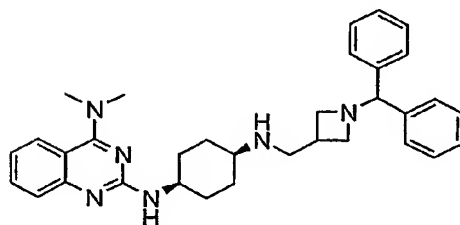
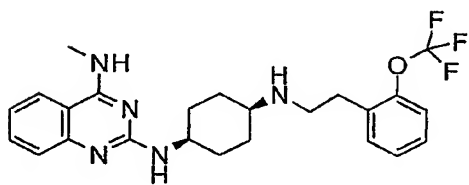
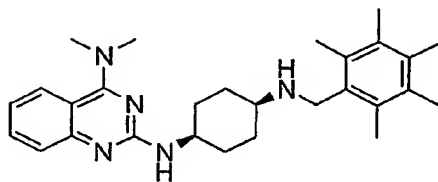
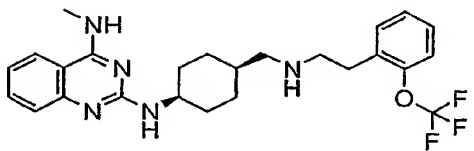
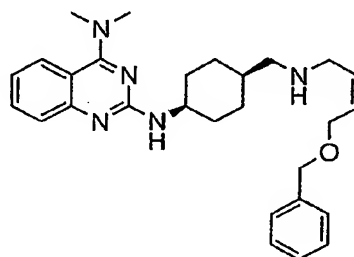
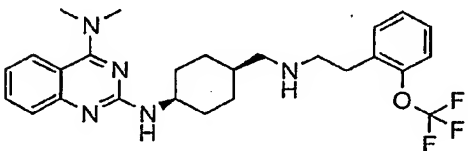
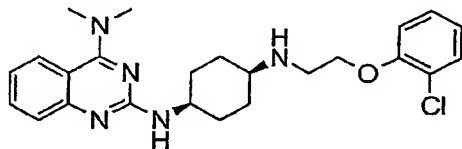
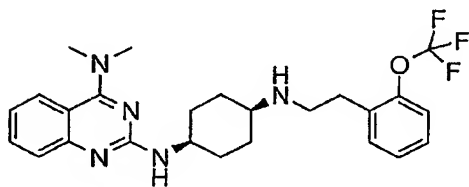
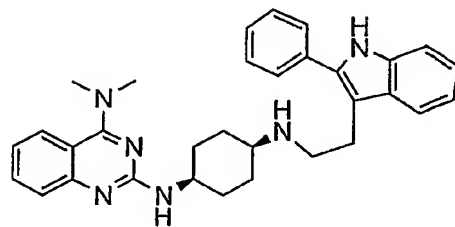
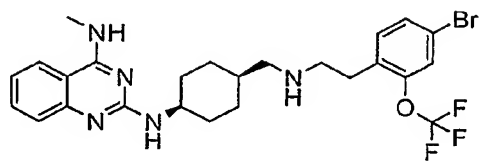
,

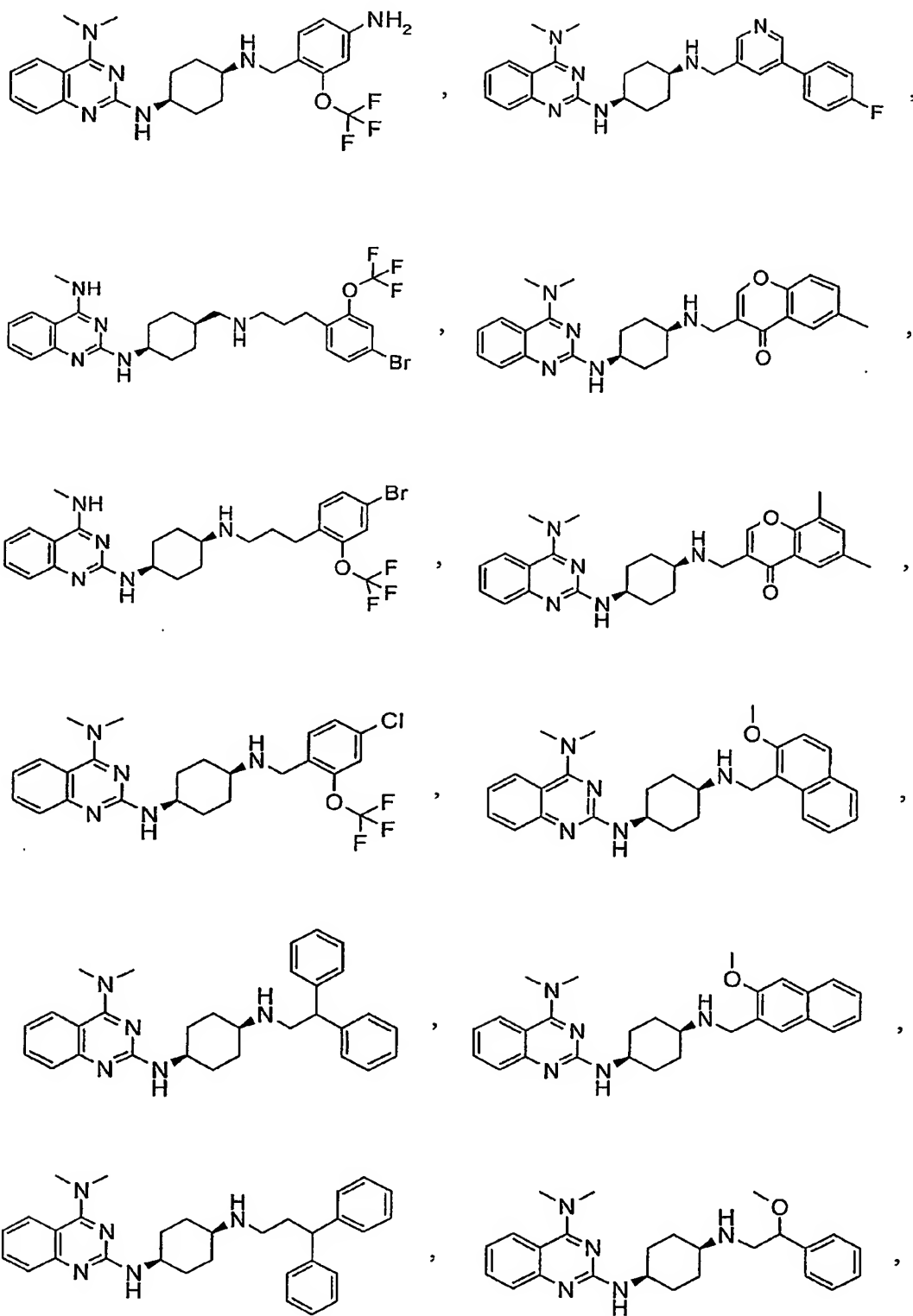


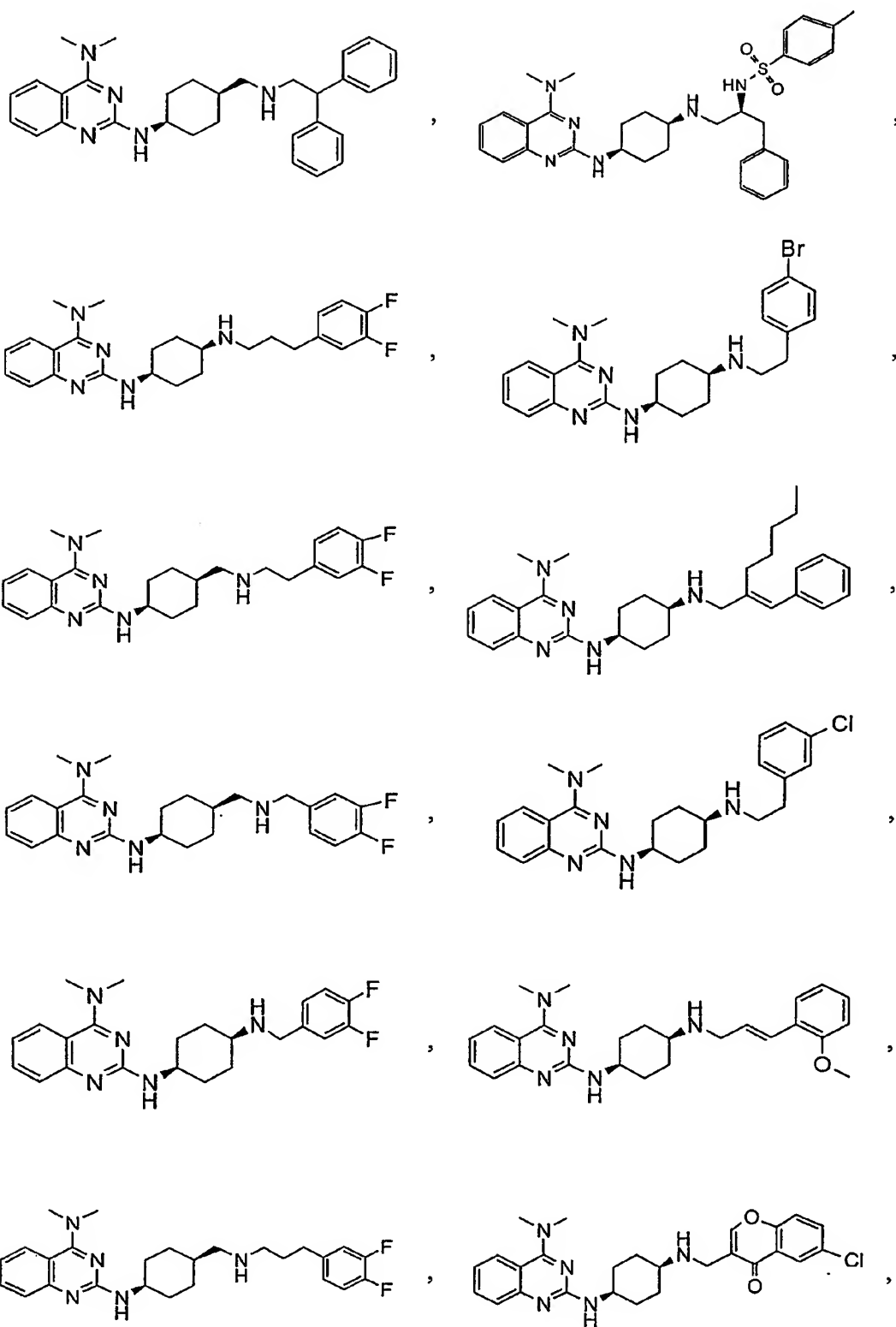
,

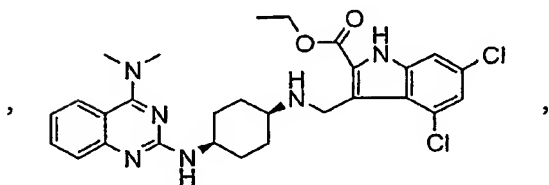
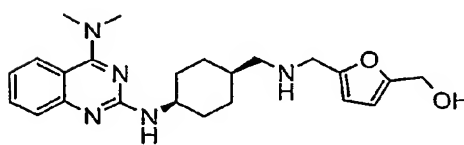
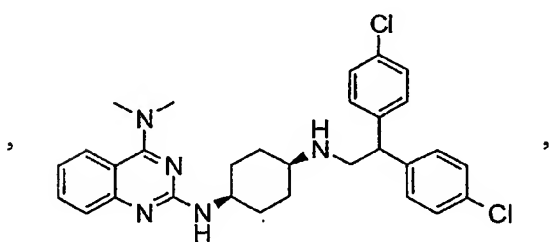
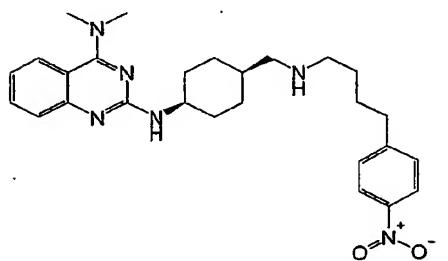
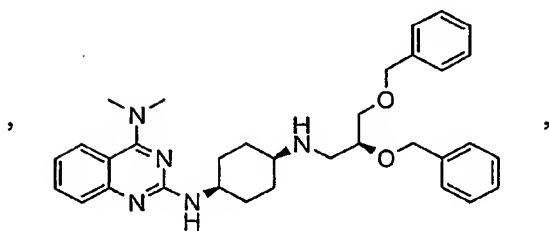
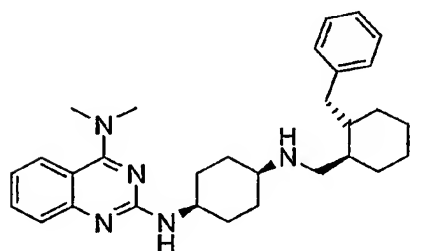
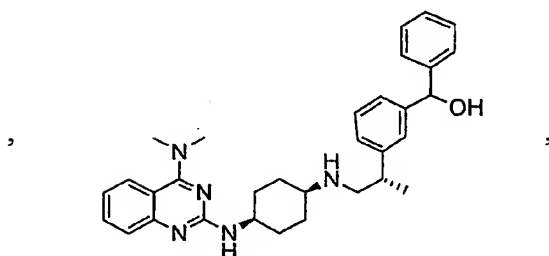
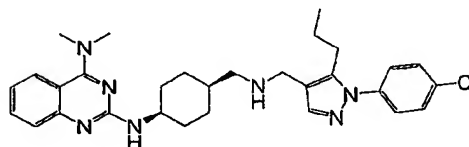
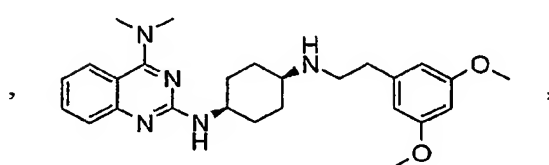
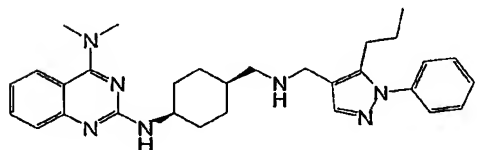
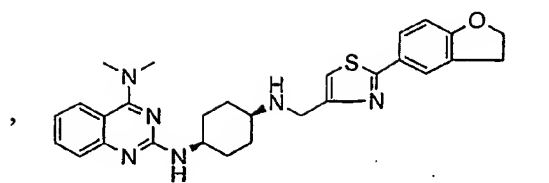
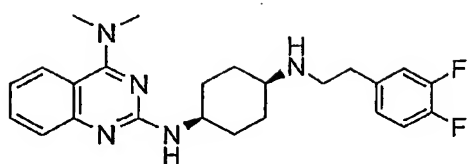


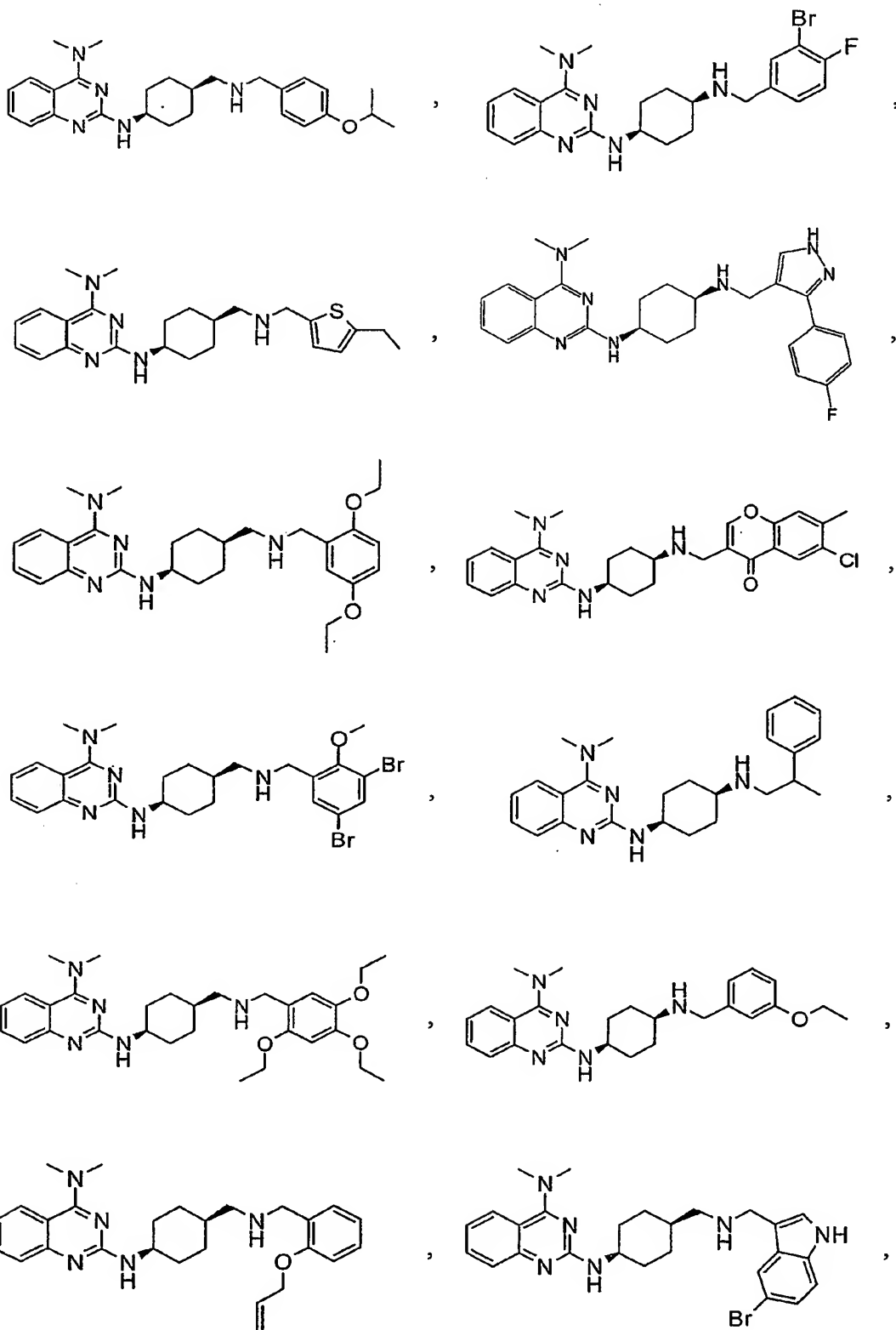


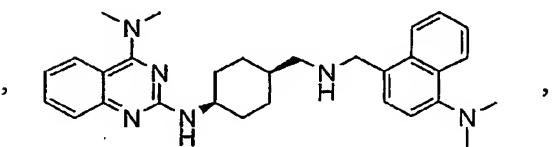
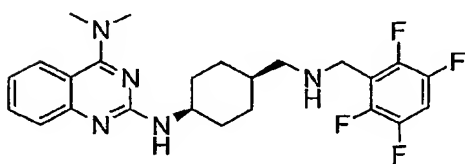
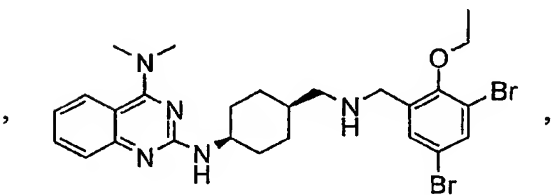
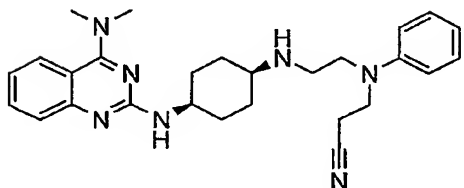
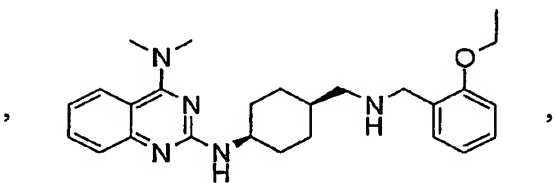
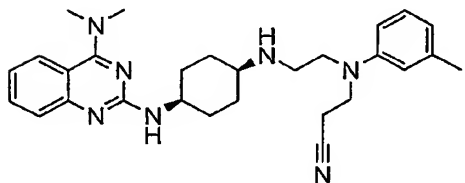
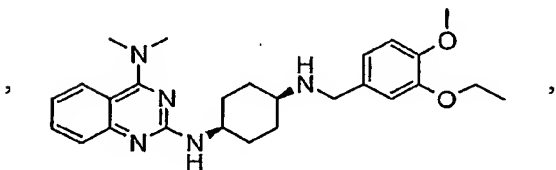
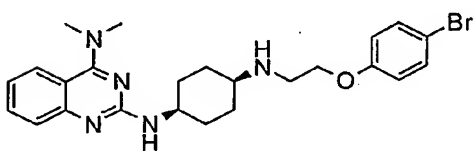
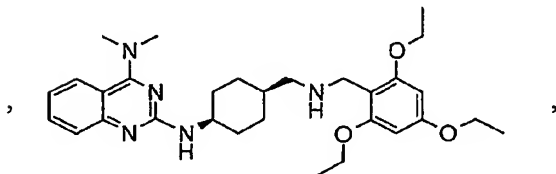
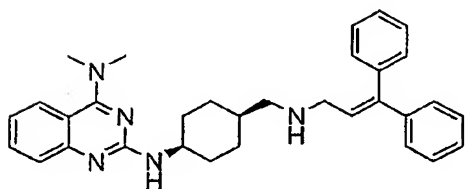
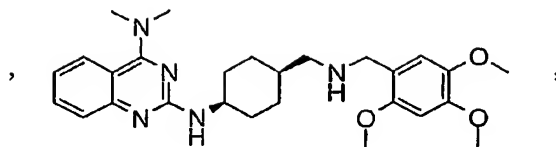
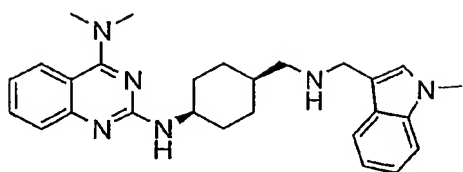


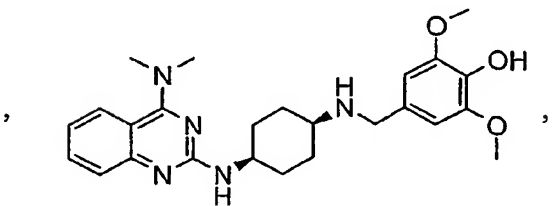
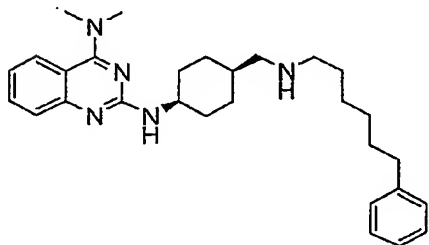
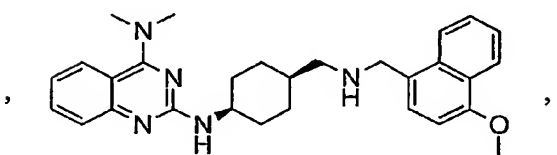
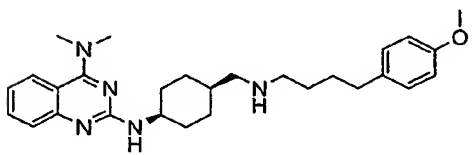
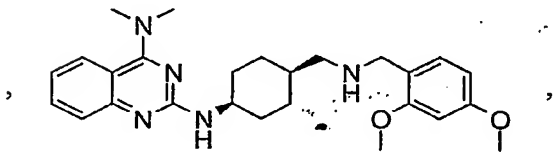
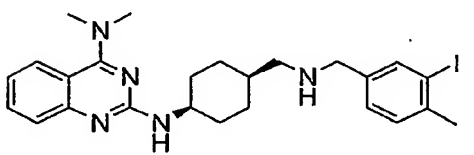
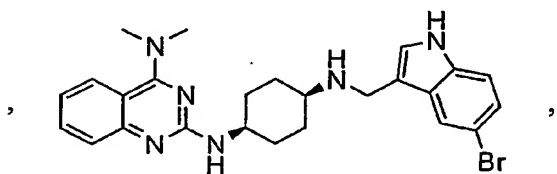
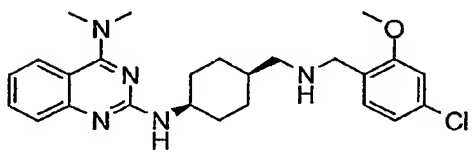
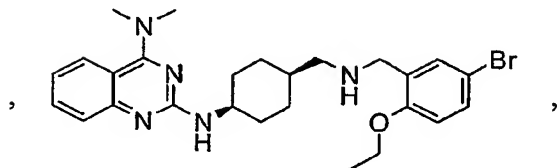
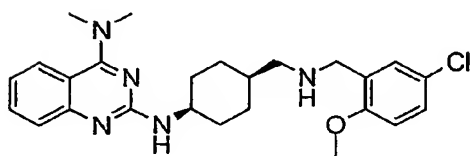
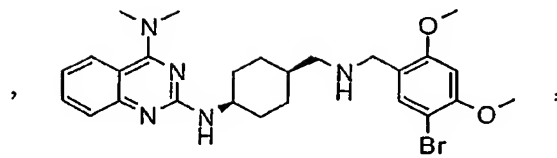
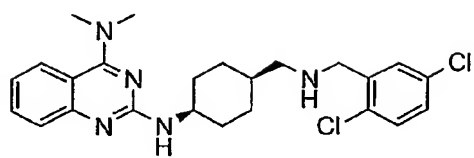


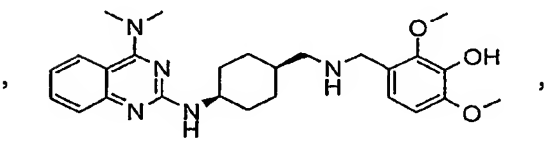
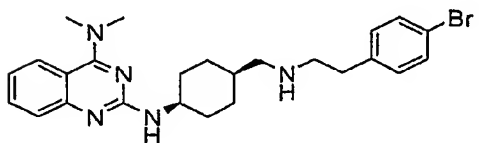
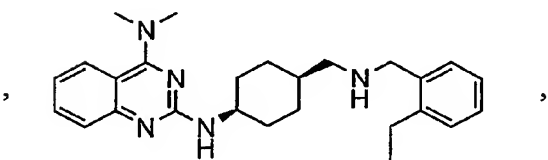
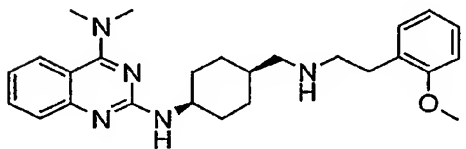
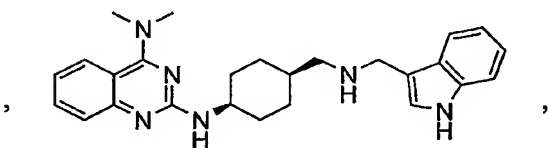
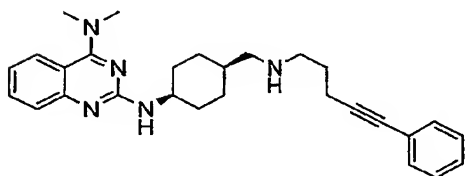
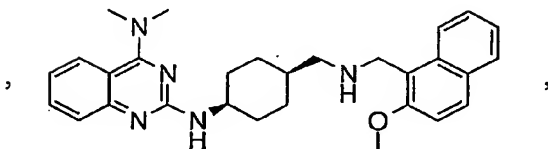
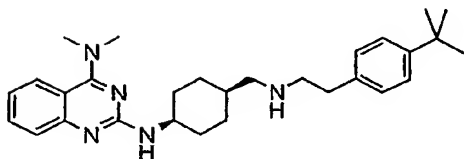
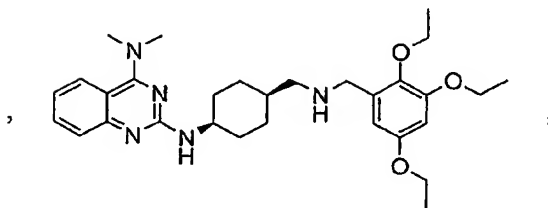
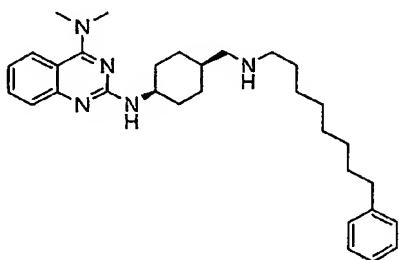
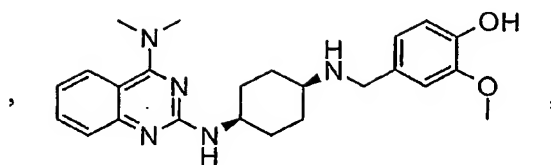
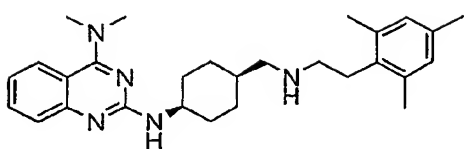


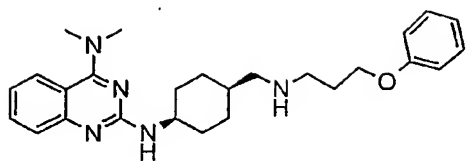




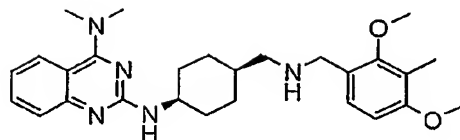




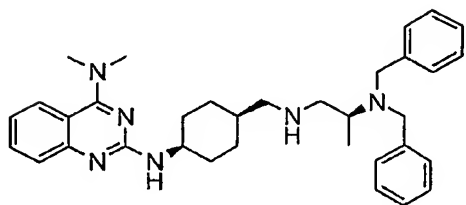




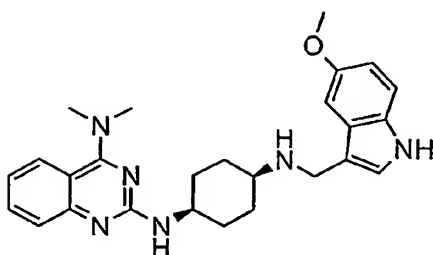
,



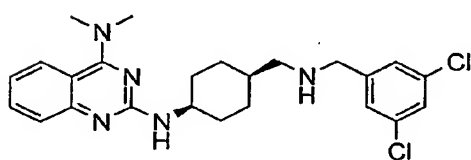
,



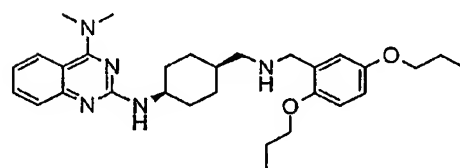
,



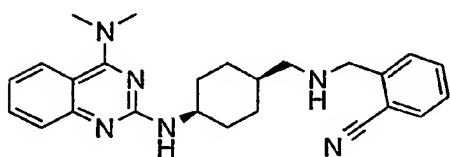
,



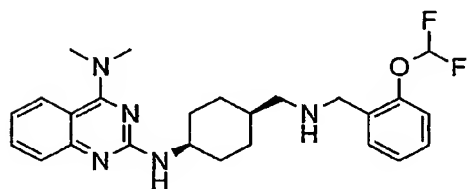
,



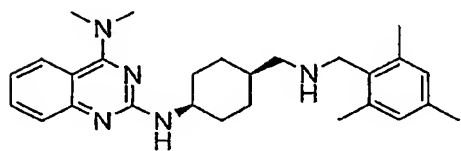
,



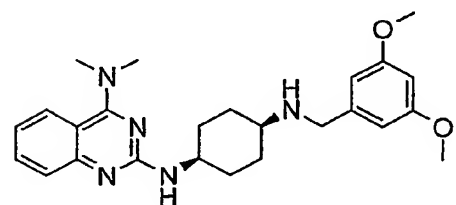
,



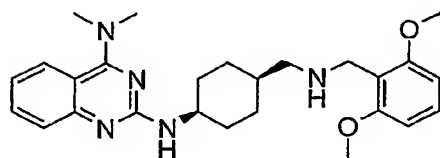
,



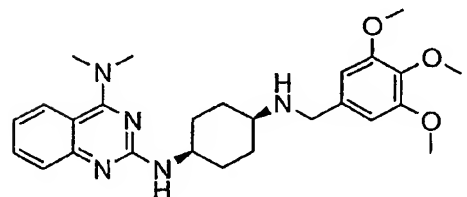
,



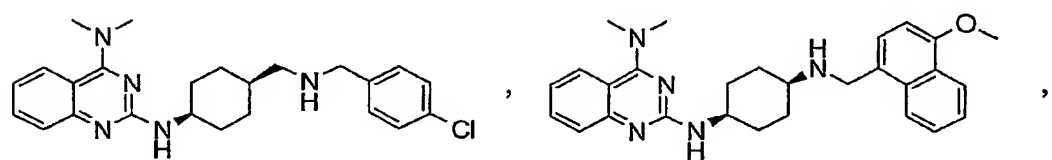
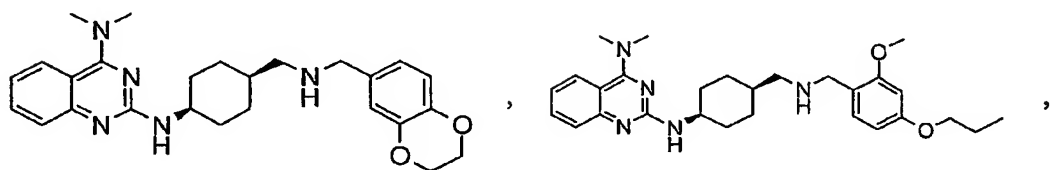
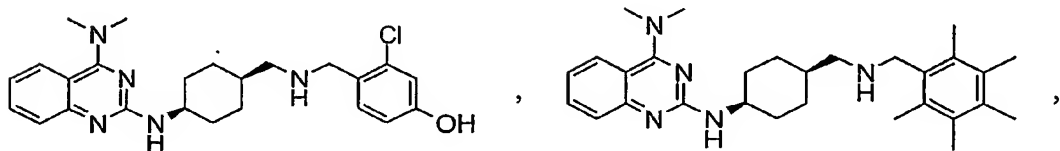
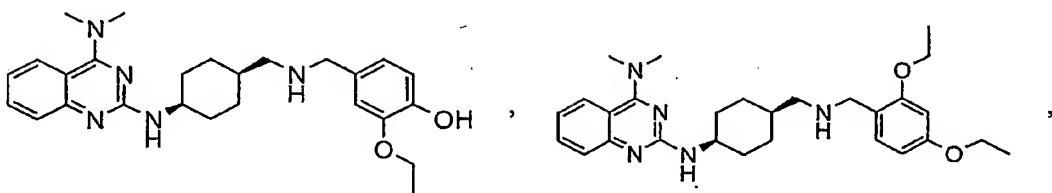
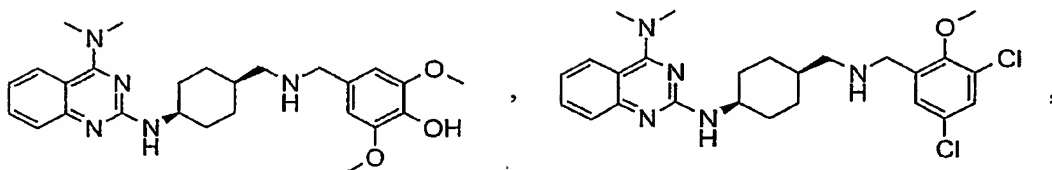
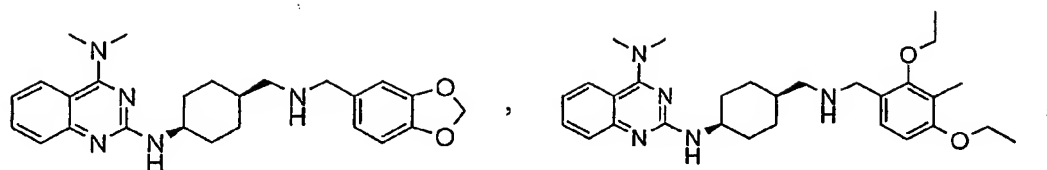
,

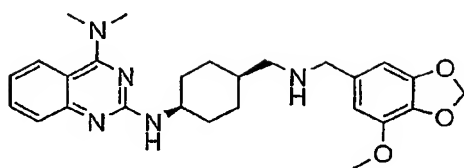


,

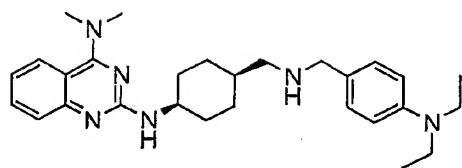


,

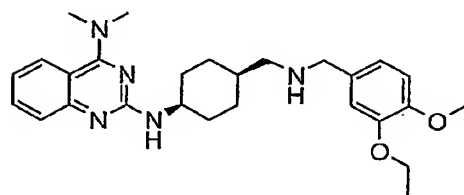




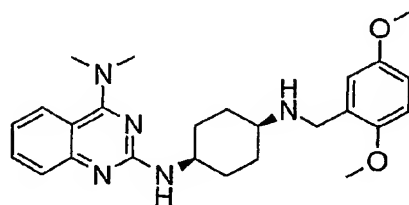
,



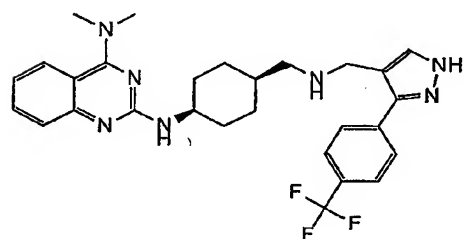
,



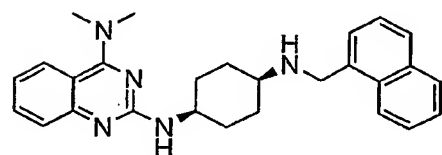
,



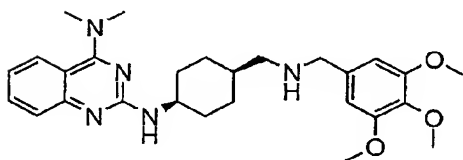
,



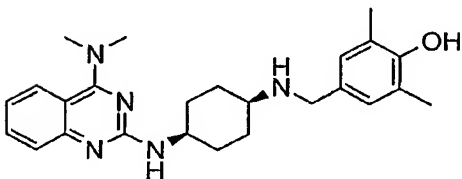
,



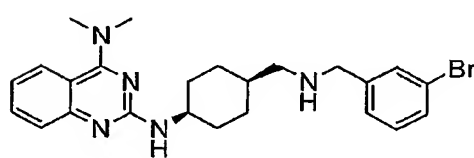
,



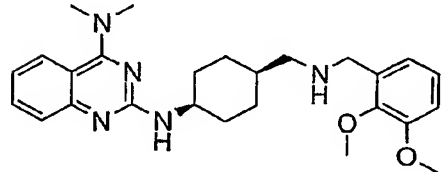
,



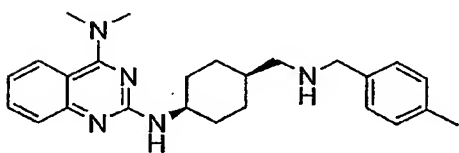
,



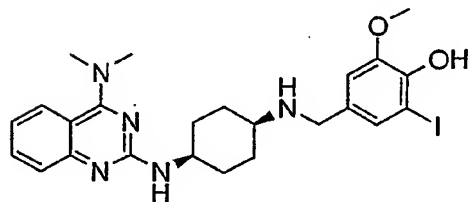
,



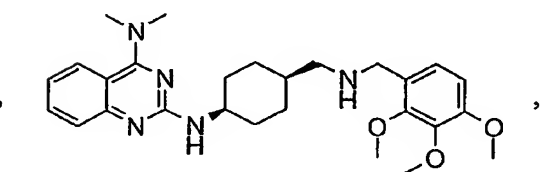
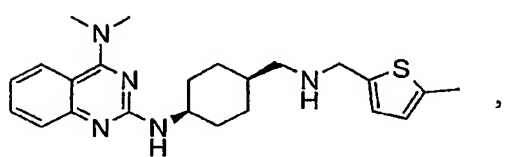
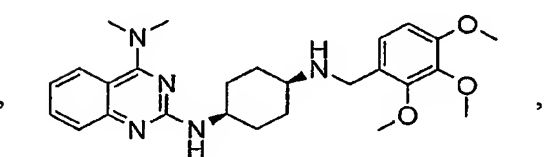
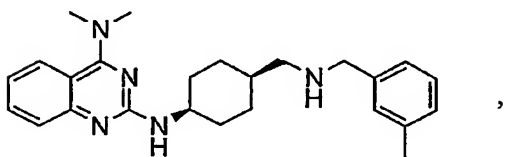
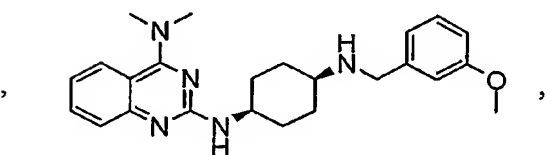
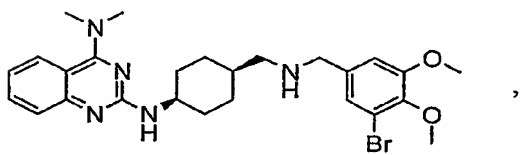
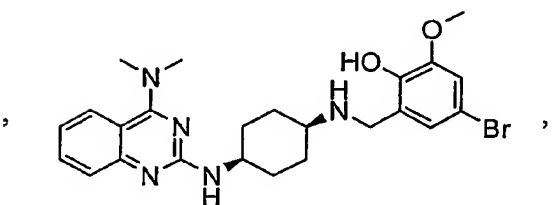
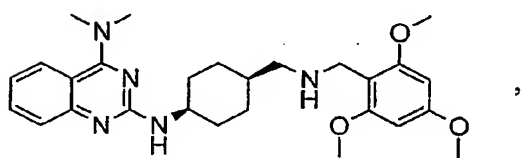
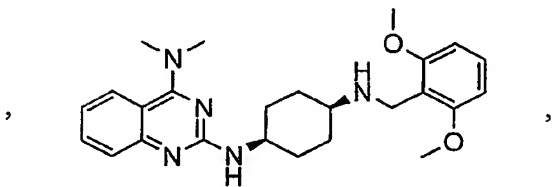
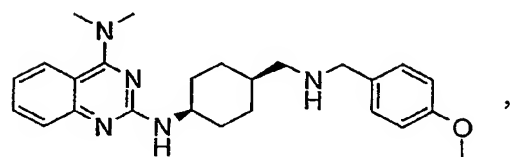
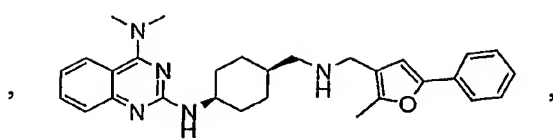
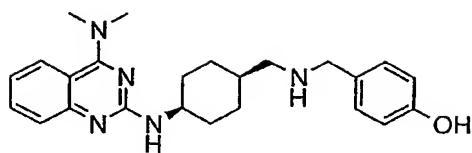
,

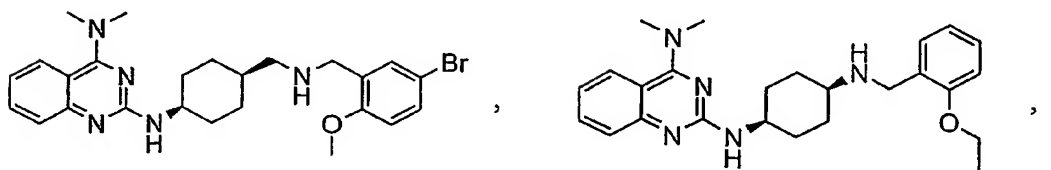
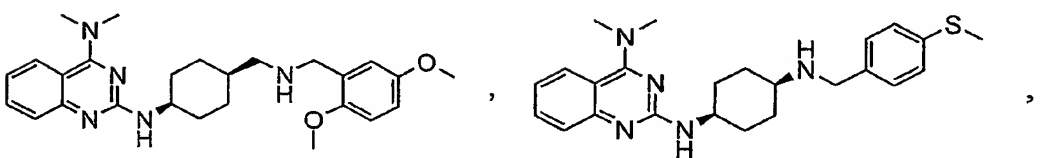
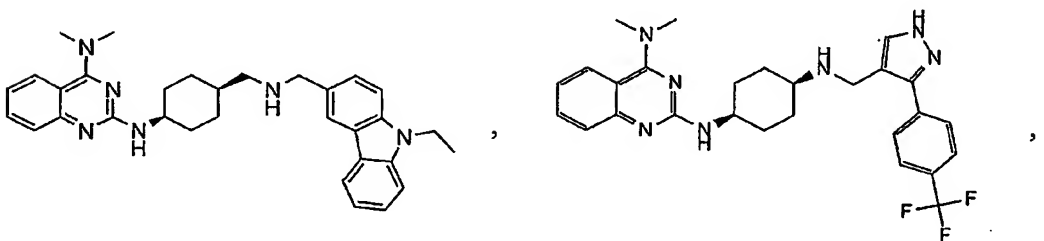
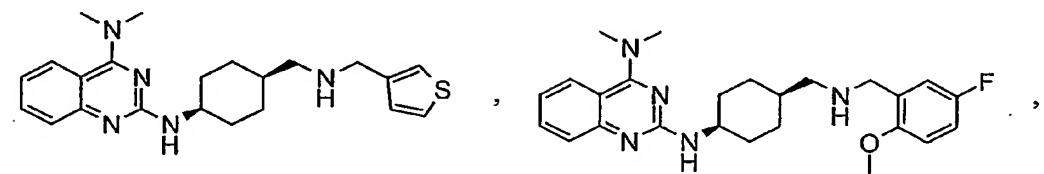
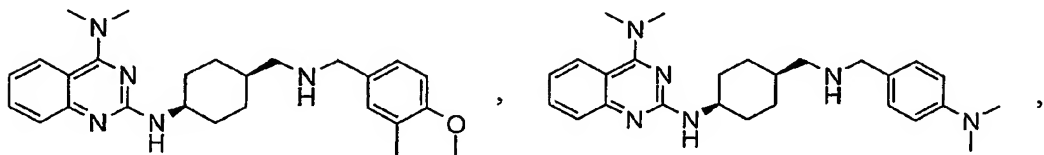
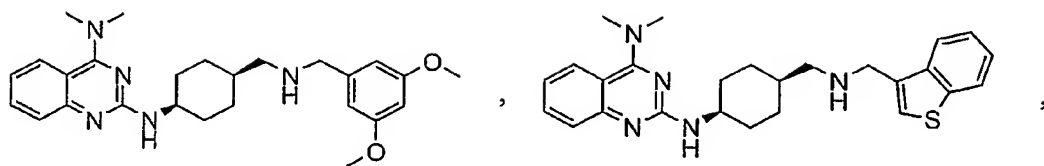


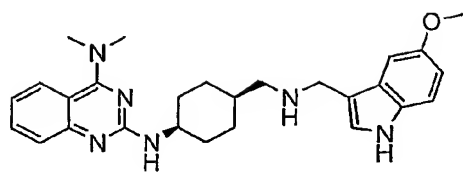
,



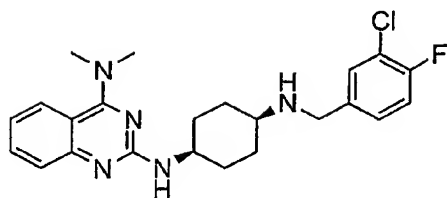
,



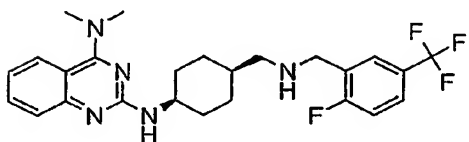




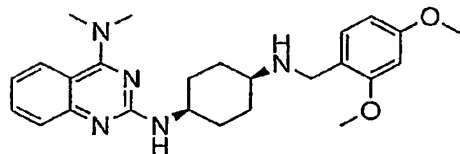
,



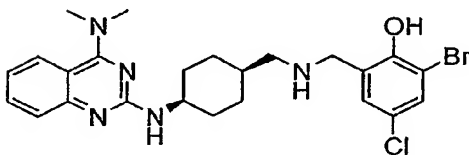
,



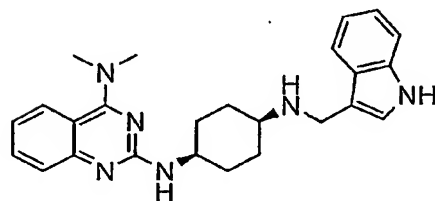
,



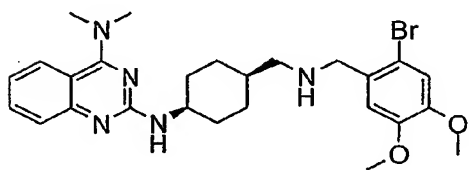
,



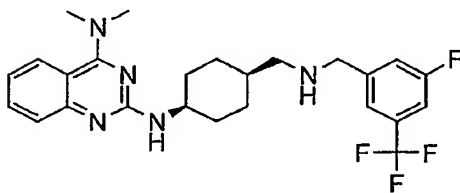
,



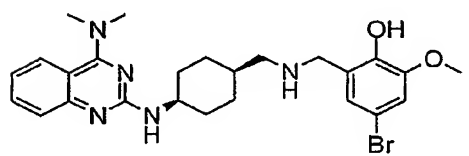
,



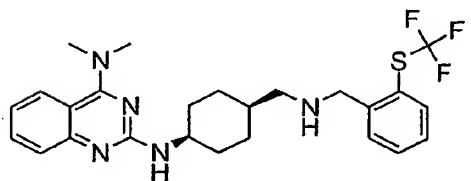
,



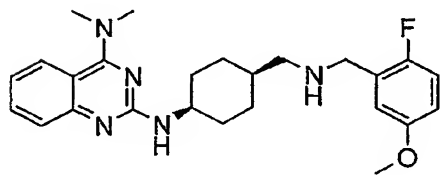
,



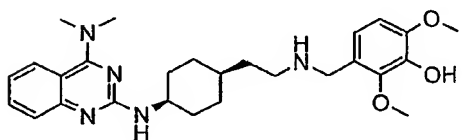
,



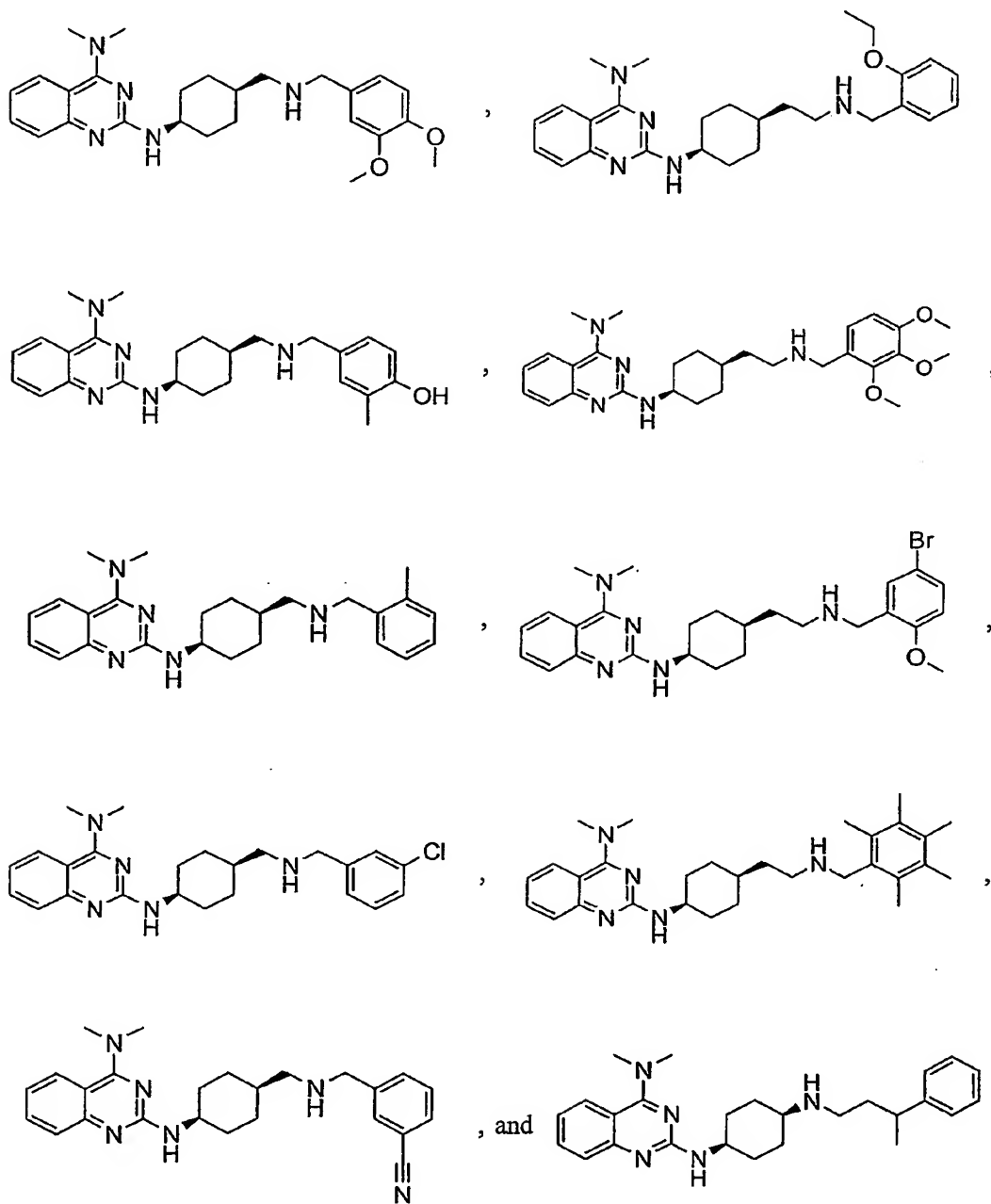
,



,



,



; or, in case of, a salt thereof.

14. A compound according to claim 1, wherein Q is Formula II;

R₁ represents

(i) C₁-C₁₆ alkyl,

C₁-C₁₆ alkyl substituted by substituent(s) independently selected from

•halogen,

•carbocyclyl,

•carbocyclic aryl,

•carbocyclic aryl substituted by substituent(s) independently selected from

••halogen,

••nitro,

••C₁-C₃ alkyl,

••halogenated C₁-C₃ alkyl,

••C₁-C₃ alkoxy,

••halogenated C₁-C₃ alkoxy,

(ii) C₂-C₃ alkenyl,

C₂-C₃ alkenyl substituted by carbocyclic aryl,

(iii) carbocyclic aryl,

carbocyclic aryl substituted by substituent(s) independently selected from

•halogen,

•cyano,

•nitro,

•C₁-C₅ alkyl,

•C₁-C₅ alkyl substituted by substituent(s) independently selected from

••halogen,

••oxo,

•C₂-C₃ alkenyl,

•C₁-C₄ alkoxy,

•C₁-C₄ alkoxy substituted by substituent(s) independently selected from

••halogen,

••heterocyclyl,

••halogenated heterocyclyl,

•carbocyclic aryloxy,

- carbocyclic aryloxy substituted by substituent(s) independently selected from

- halogen,

- nitro,

- heterocyclyloxy,

- heterocyclyloxy substituted by substituent(s) independently selected from

- halogen,

- C₁-C₃ alkyl,

- halogenated C₁-C₃ alkyl,

- C₁-C₃ alkoxycarbonyl,

- mono- or di-C₁-C₄ alkylamino,

- C₁-C₃ alkylcarbonylamino,

- carbocyclic aryl diazo,

- carbocyclic aryl diazo substituted by mono- or di- C₁-C₃ alkylamino,

- C₁-C₃ alkylsulfonyl,

- carbocyclic aryl,

- (iv) heterocyclyl,

- or heterocyclyl substituted by substituent(s) independently selected from

- halogen,

- C₁-C₃ alkyl,

- C₁-C₃ alkyl substituted by substituent(s) independently selected from

- halogen,

- oxo,

- carbocyclic arylcarbonylamino,

- halogenated carbocyclic arylcarbonylamino,

- heterocyclyl,

- heterocyclyl substituted by substituent(s) independently selected from

- halogen,

- C₁-C₃ alkyl,

- halogenated C₁-C₃ alkyl,

- C₁-C₃ alkoxy,

- C₁-C₃ alkylcarbonylamino,

- carbocyclic arylsulfonyl,

- C₁-C₃ alkoxy carbonyl,
- carbocyclic aryl,
- halogenated carbocyclic aryl,
- heterocyclyl,
- heterocyclyl substituted by substituent(s) independently selected from
 - halogen,
 - C₁-C₃ alkyl,
 - halogenated C₁-C₃ alkyl;

Y is -S(O)₂-;

wherein carbocyclic aryl is phenyl, biphenyl, or naphthyl;

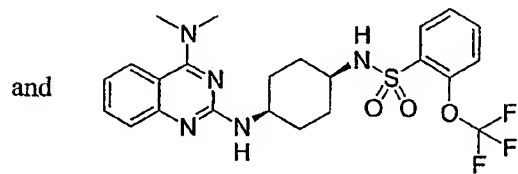
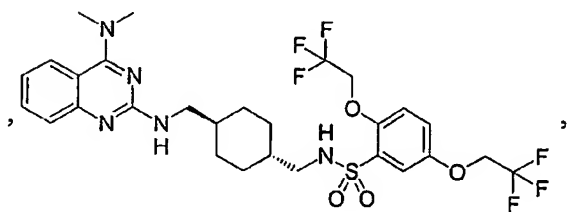
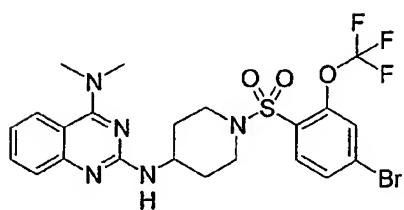
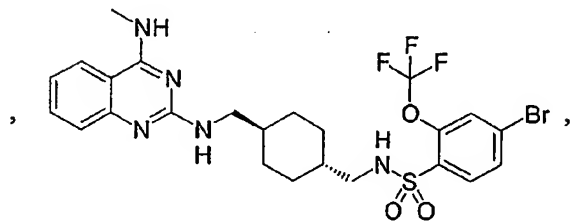
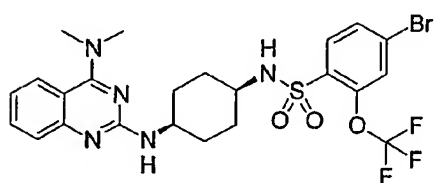
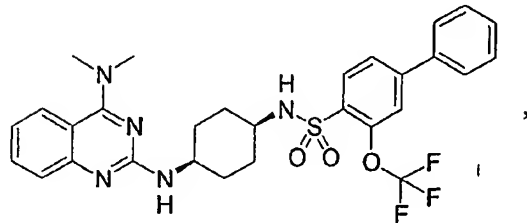
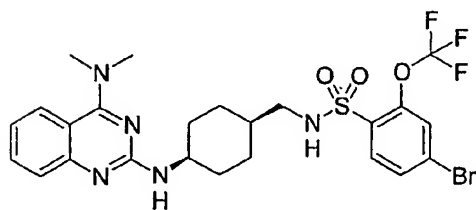
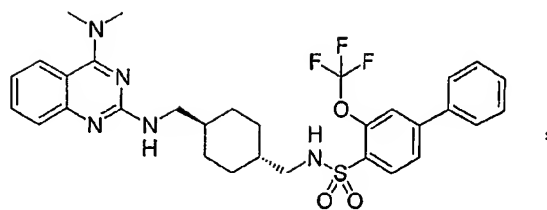
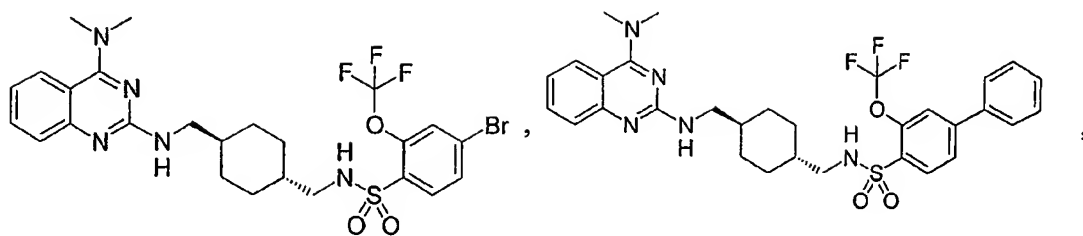
carbocyclyl is 7,7-dimethyl-2-oxo-bicyclo[2.2.1]heptyl;

heterocyclyl is 1,2,3,4-tetrahydro-isoquinolyl, 1,2,3-thiadiazolyl, 1*H*-pyrrolyl, benzo[2,1,3]oxadiazolyl, benzo[b]thienyl, furyl, imidazolyl, isoxazolyl, pyrazolyl, pyridyl, quinolyl, thiazolyl, or thienyl;

halogen is fluoro, chloro, bromo, or iodo;

or a salt thereof.

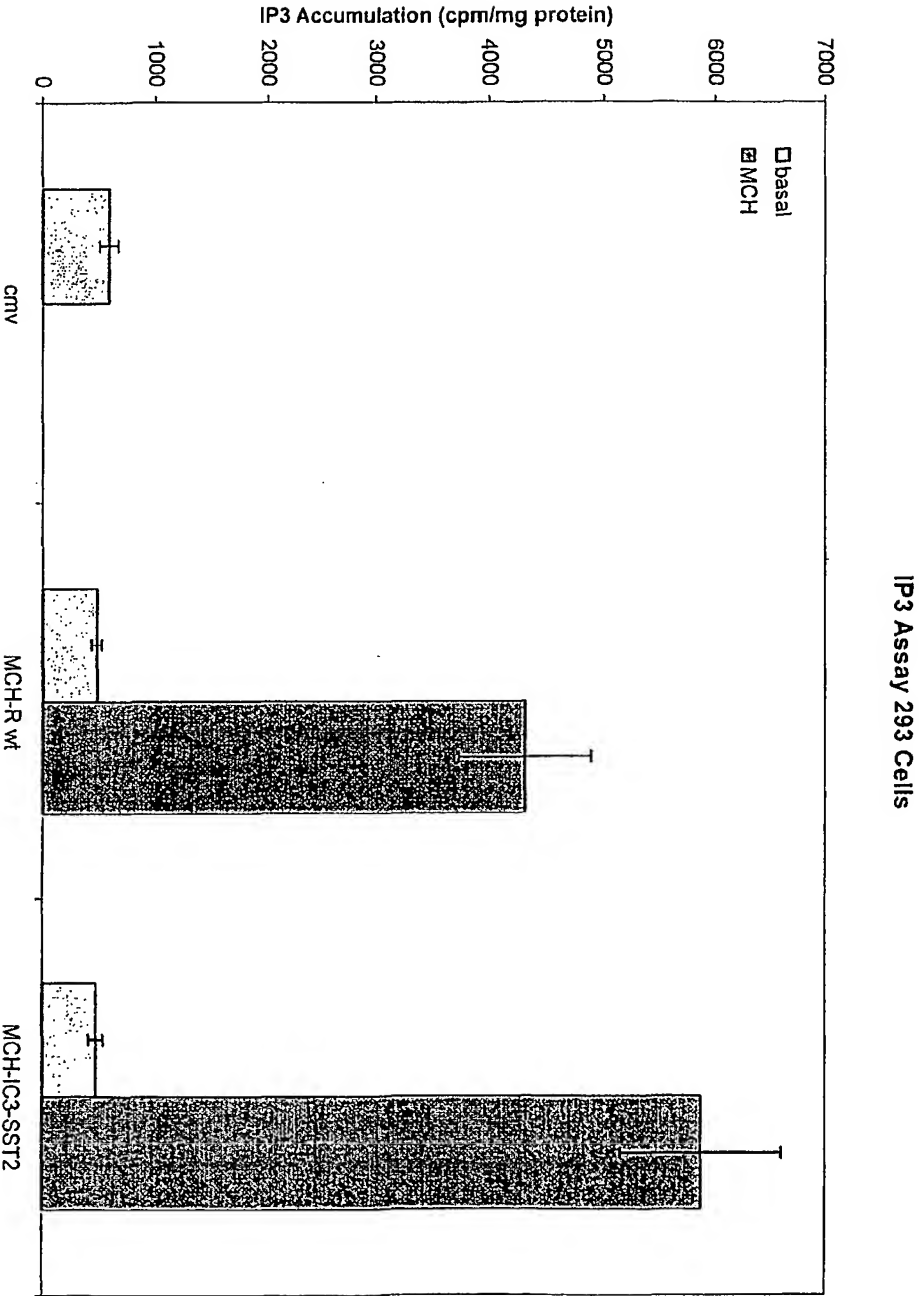
15. A compound according to claim 14 of Formua I selected from the group consisting of



; or, in case of, a salt thereof.

16. A compound according to claim 1, wherein Q is Formula II;
R₁ is selected from H, -CO₂Bu, or -CO₂Bn (Bn is a benzyl group);
R₂ is methylamino or dimethylamino;
L is selected from Formula XX - XXII;
Y is a single bond;
or a salt thereof.
17. A method for modulating the G-protein receptor, SLC-1, comprising the step of contacting said SLC-1 with a MCH receptor antagonist.
18. A method for modulating the G-protein receptor, SLC-1, comprising the step of contacting said SLC-1 with a compound of claims 1-16.
19. The method of prophylaxis or treatment of obesity, obesity related disorders, anxiety, or depression in mammals in need of such treatment comprising administering to the mammal a therapeutically effective amount of a compound having the composition of any of claims 1-16.
20. A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a therapeutically effective amount of a compound having the composition of any of claims 1-16.

Fig. 1



SEQUENCE LISTING

<110> Arena Pharmaceuticals, Inc.

<120> MCH Receptor Antagonists

<130> AREN-0238

<160> 12

<170> PatentIn version 3.0

<210> 1

<211> 30

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 1

gtgaagcttg cctctggtgc ctgcaggagg

30

<210> 2

<211> 31

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 2

gcagaattcc cggtagcggtg ttgtggtgcc c

31

<210> 3

<211> 24

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 3

catgagctgg tggatcatga aggg

24

<210> 4

<211> 24

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 4

atgaagggca tgcccaggag aaag

24

<210> 5

<211> 1349

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 5

atggacctgg aagcctcgct gctgccact ggtcccaatg ccagcaacac ctctgatggc

60

cccgataacc tcaacttcggc aggatcacct cctcgcacgg ggagcatctc ctacatcaac

120

```

atcatcatgc cttcgggtgtt cggcaccatc tgcctcctgg gcatcatcgg gaactccaag 180
gtcatcttcg cggtcgtgaa gaagtccaag ctgcaactgg gcaacaacgt ccccgacatc 240
ttcatcatca acctctcggg agtagatctc ctctttctcc tgggcatgcc cttcatgac 300
caccagctca tgggcaatgg ggtgtggcac tttggggaga ccatgtgcac cctcatcacg 360
gccatggatg ccaatagtca gttcaccagc acctacatcc tgaccgccat ggccattgac 420
cgctacctgg ccactgtcca ccccatctct tccacgaagt tccggaagcc ctctgtggcc 480
accctgggta tctgcctcct gtggggccctc tccttcatca gcatcaccoc tgtgtggctg 540
tatgccagac tcatcccctt cccaggaggt gcagtgggct gcggcatacg cctgccaac 600
ccagacactg acctctactg gttcacctg taccagtttt tcctggcctt tgccctgct 660
tttgtgttca tcacagccgc atactgagg atcctgcaga aggtgaagtc ctctggaatc 720
cgagtgggct cctctaagag gaagaagtct gagaagaagg tcaccgcac agccatcgcc 780
atctgtctgg tcttctttgt gtgctgggca ccctactatg tgctacagct gaccagttg 840
tccatcagcc gcccgacct cacctttgtc tacttataca atgcggccat cagcttgggc 900
tatgccaaac gctgcctcaa cccctttgtg tacatcgtgc tctgtgagac gttccgcaaa 960
cgcttgggtc tgtcggtgaa gcctgcagcc caggggcagc ttcgcgctgt cagcaacgct 1020
cagacggctg acgaggagag gacagaaagc aaaggcacct gatacttccc ctgccaccct 1080
gcacacctcc aagtcagggc accacaacac gccaccggga gagatgctga gaaaaacca 1140
agaccgctcg ggaaatgcag gaaggccggg ttgtgagggg ttgttgcaat gaaataaata 1200
cattccatgg gctcacacgt tgctggggag gcctggagtc aggtttgggg ttttcagata 1260
tcagaaatcc cttgggggag caggatgaga cctttggata gaacagaagc tgagcaagag 1320
aacatgttgg tttggataac cggttgcac 1349

```

<210> 6

<211> 446

<212> PRT

<213> Homo Sapien

<220>

<223> Novel Sequence

<400> 6

```

Met Asp Leu Glu Ala Ser Leu Leu Pro Thr Gly Pro Asn Ala Ser Asn
 1           5           10           15

Thr Ser Asp Gly Pro Asp Asn Leu Thr Ser Ala Gly Ser Pro Pro Arg
          20           25           30

Thr Gly Ser Ile Ser Tyr Ile Asn Ile Ile Met Pro Ser Val Phe Gly
      35           40           45

Thr Ile Cys Leu Leu Gly Ile Ile Gly Asn Ser Thr Val Ile Phe Ala
 50           55           60

Val Val Lys Lys Ser Lys Leu His Trp Cys Asn Asn Val Pro Asp Ile
 65           70           75           80

Phe Ile Ile Asn Leu Ser Val Val Asp Leu Leu Phe Leu Leu Gly Met
          85           90           95

Pro Phe Met Ile His Gln Leu Met Gly Asn Gly Val Trp His Phe Gly
      100           105           110

Glu Thr Met Cys Thr Leu Ile Thr Ala Met Asp Ala Asn Ser Gln Phe
      115           120           125

Thr Ser Thr Tyr Ile Leu Thr Ala Met Ala Ile Asp Arg Tyr Leu Ala
      130           135           140

Thr Val His Pro Ile Ser Ser Thr Lys Phe Arg Lys Pro Ser Val Ala
      145           150           155           160

Thr Leu Val Ile Cys Leu Leu Trp Ala Leu Ser Phe Ile Ser Ile Thr
          165           170           175

Pro Val Trp Leu Tyr Ala Arg Leu Ile Pro Phe Pro Gly Gly Ala Val
      180           185           190

Gly Cys Gly Ile Arg Leu Pro Asn Pro Asp Thr Asp Leu Tyr Trp Phe
      195           200           205

Thr Leu Tyr Gln Phe Phe Leu Ala Phe Ala Leu Pro Phe Val Val Ile
      210           215           220

Thr Ala Ala Tyr Val Arg Ile Leu Gln Lys Val Lys Ser Ser Gly Ile
      225           230           235           240

Arg Val Gly Ser Ser Lys Arg Lys Lys Ser Glu Lys Lys Val Thr Arg
          245           250           255

Thr Ala Ile Ala Ile Cys Leu Val Phe Phe Val Cys Trp Ala Pro Tyr
      260           265           270

Tyr Val Leu Gln Leu Thr Gln Leu Ser Ile Ser Arg Pro Thr Leu Thr
      275           280           285

Phe Val Tyr Leu Tyr Asn Ala Ala Ile Ser Leu Gly Tyr Ala Asn Ser

```

290		295		300
Cys Leu Asn Pro Phe Val Tyr Ile Val Leu Cys Glu Thr Phe Arg Lys				
305		310		315 320
Arg Leu Val Leu Ser Val Lys Pro Ala Ala Gln Gly Gln Leu Arg Ala				
	325		330	335
Val Ser Asn Ala Gln Thr Ala Asp Glu Glu Arg Thr Glu Ser Lys Gly				
	340		345	350
Thr Tyr Phe Pro Cys His Pro Ala His Leu Gln Val Arg Ala Pro Gln				
	355		360	365
His Ala Thr Gly Arg Asp Ala Glu Lys Asn Pro Arg Pro Leu Gly Lys				
	370		375	380
Cys Arg Lys Ala Gly Leu Gly Val Val Ala Met Lys Ile His Ser Met				
385		390		395 400
Gly Ser His Val Ala Gly Glu Ala Trp Ser Gln Val Trp Gly Phe Gln				
	405		410	415
Ile Ser Glu Ile Pro Trp Gly Ser Arg Met Arg Pro Leu Asp Arg Thr				
	420		425	430
Glu Ala Glu Gln Glu Asn Met Leu Val Trp Ile Thr Gly Cys				
	435		440	445

<210> 7
 <211> 70

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 7
 gatcctgcag aaggtgaagt cctctggaat ccgagtgggc tcctctaaga ggaagaagtc 60
 tgagaagaag 70

<210> 8

<211> 71

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 8

gtgaccttct tctcagactt cttcctctta gaggagccca ctoggattcc agaggacttc 60

accttctgca g 71

<210> 9

<211> 30

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 9

gtgaagcttg cccgggcagg atggacctgg 30

<210> 10

<211> 24

<212> DNA

<213> Artificial

<220>

<223> Novel Sequence

<400> 10

atctagaggt gcctttgctt tctg 24

<210> 11

<211> 2133

<212> DNA

<213> Homo Sapien

<400> 11

atggacctgg aagcctcgct gctgcccact ggtcccaatg ccagcaaac ctctgatggc 60

cccgataacc tcaacttcggc aggatcacct cctcgcacgg ggagcatctc ctacatcaac 120

atcatcatgc cttcgggtgtt cggcaccatc tgccctcctgg gcatcatcgg gaactccacg 180
gtcatcttcg cggtcgtgaa gaagtccaag ctgcactggg gcaacaacgt ccccgacatc 240
ttcatcatca acctctcggg agtagatctc ctctttctcc tgggcatgcc cttcatgacg 300
caccagctca tgggcaatgg ggtgtggcac tttggggaga ccatgtgcac cctcatcacg 360
gccatggatg ccaatagtca gttcaccagc acctacatcc tgaccgccat ggccattgac 420
cgctacctgg ccactgtcca ccccatctct tccacgaagt tccggaagcc ctctgtggcc 480
accctgggtga tctgcctcct gtggggccctc tccttcatca gcatcacccc tgtgtggctg 540
tatgccagac tcatcccctt cccaggagggt gcagtgggct gcggcatacg cctgcccac 600
ccagacactg acctctactg gttcaccctg taccagtttt tccctggcctt tgccctgcct 660
tttgtgggtca tcacagccgc atacgtgagg atcctgcagc gcatgacgtc ctcatgtggc 720
cccgccctcc agcgagcat ccggctgcgg acaaagaggg tgaccgcac agccatcgcc 780
atctgtctgg tcttctttgt gtgctgggca cctactatg tgctacagct gaccagttg 840
tccatcagcc gcccgacct cacctttgtc tacttataca atgcggccat cagcttgggc 900
tatgccaaca gctgcctcaa cccctttgtg tacatcgtgc tctgtgagac gttccgcaa 960
cgcttggtcc tgtcgggtgaa gcctgcagcc caggggcagc ttcgcgtgt cagcaacgct 1020
cagacggctg acgaggagag gacagaaagc aaaggcacct ctagaatggg ctgcacactg 1080
agcgctgagg acaaggcggc cgtggagcgc agcaagatga tcgaccgca cctccgggag 1140
gacggagaga aggcagcgc cgaggtaag ctgctgctgc tgggtgctgg tgaatccggg 1200
aagagcacia ttgtgaagca gatgaaaatt atccacgagg ctggctactc agaggaagag 1260
tgtaagcagt acaaagcagt ggtctacagc aacaccatcc agtccatcat tgccatcatt 1320
agagccatgg ggagattgaa aatcgacttt ggagacgctg ctgctgcgga tgatgctgc 1380
caactcttcg tgcttgctgg ggctgcagag gaaggcttta tgaccgcgga gctcgccggc 1440
gtcataaaga gactgtggaa ggacagcggg gtgcaagcct gcttcaacag atcccgggag 1500
taccagctga acgattcggc ggcgtactac ctgaatgact tggacagaat agcacaacca 1560
aattacatcc caaccagca ggatgttctc agaactagag tgaaaacgac ggggaattgtg 1620
gaaacccact ttactttcaa agatcttcat tttaaaatgt ttgacgtggg aggccagaga 1680
tcagagcggg agaagtggat tcaactgcttt gaaggcgtga ctgccatcat cttctgtgtg 1740
gccctgagtg actatgacct ggttcttgct gaggatgaag aaatgaaccg gatgcatgaa 1800
agcatgaagc tgttcgatag catatgtaac aacaagtggg ttacggacac atccatcatc 1860

cttttcctga acaagaagga cctcttcgaa gagaagatca aaaagagtcc cctcacgata 1920
 tgctatccag aatatgcagg ctcaaacaca tatgaagagg cggtgcgta tatccagtgt 1980
 cagtttgaag acctcaataa aaggaaggac acaaaggaaa ttacaccca cttcacttgc 2040
 gccacggata cgaagaatgt gcagtttgtg ttcgatgctg taacggacgt catcataaag 2100
 aataacctaa aagactgtgg tctcttctaa tct 2133

<210> 12
 <211> 709
 <212> PRT
 <213> Homo Sapien

<400> 12
 Met Asp Leu Glu Ala Ser Leu Leu Pro Thr Gly Pro Asn Ala Ser Asn
 1 5 10 15
 Thr Ser Asp Gly Pro Asp Asn Leu Thr Ser Ala Gly Ser Pro Pro Arg
 20 25 30
 Thr Gly Ser Ile Ser Tyr Ile Asn Ile Ile Met Pro Ser Val Phe Gly
 35 40 45
 Thr Ile Cys Leu Leu Gly Ile Ile Gly Asn Ser Thr Val Ile Phe Ala
 50 55 60
 Val Val Lys Lys Ser Lys Leu His Trp Cys Asn Asn Val Pro Asp Ile
 65 70 75 80
 Phe Ile Ile Asn Leu Ser Val Val Asp Leu Leu Phe Leu Leu Gly Met
 85 90 95
 Pro Phe Met Ile His Gln Leu Met Gly Asn Gly Val Trp His Phe Gly
 100 105 110
 Glu Thr Met Cys Thr Leu Ile Thr Ala Met Asp Ala Asn Ser Gln Phe
 115 120 125
 Thr Ser Thr Tyr Ile Leu Thr Ala Met Ala Ile Asp Arg Tyr Leu Ala
 130 135 140
 Thr Val His Pro Ile Ser Ser Thr Lys Phe Arg Lys Pro Ser Val Ala
 145 150 155 160
 Thr Leu Val Ile Cys Leu Leu Trp Ala Leu Ser Phe Ile Ser Ile Thr
 165 170 175
 Pro Val Trp Leu Tyr Ala Arg Leu Ile Pro Phe Pro Gly Gly Ala Val
 180 185 190
 Gly Cys Gly Ile Arg Leu Pro Asn Pro Asp Thr Asp Leu Tyr Trp Phe
 195 200 205

Thr Leu Tyr Gln Phe Phe Leu Ala Phe Ala Leu Pro Phe Val Val Ile
 210 215 220
 Thr Ala Ala Tyr Val Arg Ile Leu Gln Arg Met Thr Ser Ser Val Ala
 225 230 235 240
 Pro Ala Ser Gln Arg Ser Ile Arg Leu Arg Thr Lys Arg Val Thr Arg
 245 250 255
 Thr Ala Ile Ala Ile Cys Leu Val Phe Phe Val Cys Trp Ala Pro Tyr
 260 265 270
 Tyr Val Leu Gln Leu Thr Gln Leu Ser Ile Ser Arg Pro Thr Leu Thr
 275 280 285
 Phe Val Tyr Leu Tyr Asn Ala Ala Ile Ser Leu Gly Tyr Ala Asn Ser
 290 295 300
 Cys Leu Asn Pro Phe Val Tyr Ile Val Leu Cys Glu Thr Phe Arg Lys
 305 310 315 320
 Arg Leu Val Leu Ser Val Lys Pro Ala Ala Gln Gly Gln Leu Arg Ala
 325 330 335
 Val Ser Asn Ala Gln Thr Ala Asp Glu Glu Arg Thr Glu Ser Lys Gly
 340 345 350
 Thr Ser Arg Met Gly Cys Thr Leu Ser Ala Glu Asp Lys Ala Ala Val
 355 360 365
 Glu Arg Ser Lys Met Ile Asp Arg Asn Leu Arg Glu Asp Gly Glu Lys
 370 375 380
 Ala Ala Arg Glu Val Lys Leu Leu Leu Leu Gly Ala Gly Glu Ser Gly
 385 390 395 400
 Lys Ser Thr Ile Val Lys Gln Met Lys Ile Ile His Glu Ala Gly Tyr
 405 410 415
 Ser Glu Glu Glu Cys Lys Gln Tyr Lys Ala Val Val Tyr Ser Asn Thr
 420 425 430
 Ile Gln Ser Ile Ile Ala Ile Ile Arg Ala Met Gly Arg Leu Lys Ile
 435 440 445
 Asp Phe Gly Asp Ala Ala Arg Ala Asp Asp Ala Arg Gln Leu Phe Val
 450 455 460
 Leu Ala Gly Ala Ala Glu Glu Gly Phe Met Thr Ala Glu Leu Ala Gly
 465 470 475 480
 Val Ile Lys Arg Leu Trp Lys Asp Ser Gly Val Gln Ala Cys Phe Asn
 485 490 495
 Arg Ser Arg Glu Tyr Gln Leu Asn Asp Ser Ala Ala Tyr Tyr Leu Asn
 500 505 510

Asp Leu Asp Arg Ile Ala Gln Pro Asn Tyr Ile Pro Thr Gln Gln Asp
 515 520 525
 Val Leu Arg Thr Arg Val Lys Thr Thr Gly Ile Val Glu Thr His Phe
 530 535 540
 Thr Phe Lys Asp Leu His Phe Lys Met Phe Asp Val Gly Gly Gln Arg
 545 550 555 560
 Ser Glu Arg Lys Lys Trp Ile His Cys Phe Glu Gly Val Thr Ala Ile
 565 570 575
 Ile Phe Cys Val Ala Leu Ser Asp Tyr Asp Leu Val Leu Ala Glu Asp
 580 585 590
 Glu Glu Met Asn Arg Met His Glu Ser Met Lys Leu Phe Asp Ser Ile
 595 600 605
 Cys Asn Asn Lys Trp Phe Thr Asp Thr Ser Ile Ile Leu Phe Leu Asn
 610 615 620
 Lys Lys Asp Leu Phe Glu Glu Lys Ile Lys Lys Ser Pro Leu Thr Ile
 625 630 635 640
 Cys Tyr Pro Glu Tyr Ala Gly Ser Asn Thr Tyr Glu Glu Ala Ala Ala
 645 650 655
 Tyr Ile Gln Cys Gln Phe Glu Asp Leu Asn Lys Arg Lys Asp Thr Lys
 660 665 670
 Glu Ile Tyr Thr His Phe Thr Cys Ala Thr Asp Thr Lys Asn Val Gln
 675 680 685
 Phe Val Phe Asp Ala Val Thr Asp Val Ile Ile Lys Asn Asn Leu Lys
 690 695 700
 Asp Cys Gly Leu Phe
 705